SUPPORTING INFORMATION

Discovery of First-in-Class Peptidomimetic Neurolysin Activators Possessing Enhanced Brain Penetration and Stability

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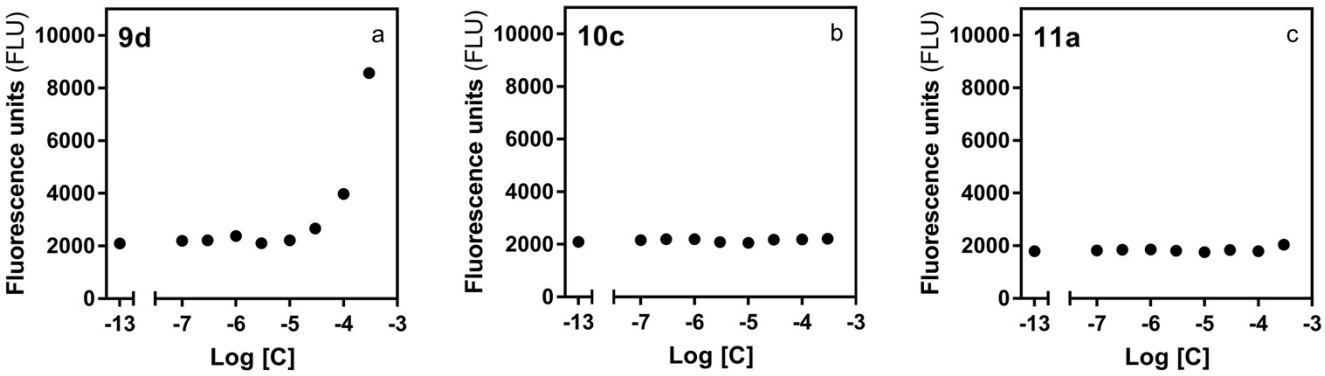
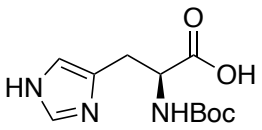
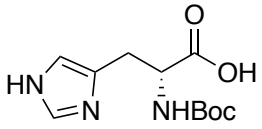


Figure S1. Interference of compounds, a) **9d**, b) **10c**, c) **11a** on the florescence intensity of the Nln hydrolytic product QFS used in the assay, a) . Compounds **10c** and **11a** provide no interference while compound **9d** interference was adjusted for in assay data analysis.

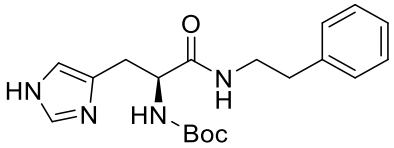
Characterization Data of Key Intermediates and Starting Material



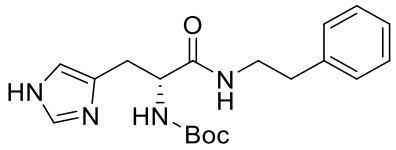
N-Boc-L-Histidine: $[\alpha]^{24}_{D}$ +24.46 (*c* 0.45, MeOH).



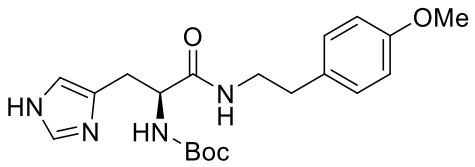
N-Boc-D-Histidine: $[\alpha]^{24}_{D}$ -24.57 (*c* 0.45, MeOH).



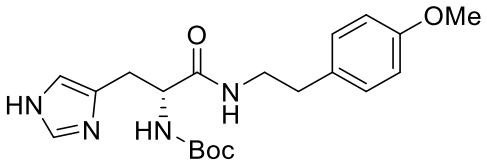
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-(phenethylamino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (84%). R_f = 0.41 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.41 (s, 9H), 2.73-2.86 (m, 4H), 3.44-3.46 (m, 2H), 4.20-4.28 (m, 1H), 6.84 (s, 1H), 7.19-7.29 (m, 5H), 7.61 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.2, 29.4, 35.1, 40.5, 54.8, 79.2, 117.0, 125.9, 128.1, 128.4, 133.2, 134.7, 139.0, 156.1, 172.8.



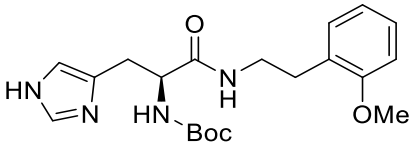
***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-(phenethylamino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (69%). R_f = 0.42 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.41 (s, 9H), 2.74-2.77 (m, 2H), 2.80-2.85 (m, 1H), 2.97-3.02 (m, 1H), 3.38-3.46 (m, 2H), 4.25 (t, *J* = 6.6 Hz, 1H), 6.84 (s, 1H), 7.18-7.30 (m, 5H), 7.61 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.25, 29.50, 35.09, 40.58, 54.87, 79.24, 116.72, 125.95, 128.12, 128.41, 134.96, 139.01, 172.86.



***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((4-methoxyphenethyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (56 %). R_f = 0.35 (MeOH : DCM = 1 : 9).^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.41 (s, 9H), 2.69 (t, *J* = 7.2 Hz, 2H), 2.80-2.86 (m, 1H), 2.98-3.03 (m, 1H), 3.34-3.40 (m, 2H), 3.76 (s, 3H), 4.25 (t, *J*= 6.8Hz, 1H), 6.84 (d, *J* = 8.4 Hz, 3H), 7.11 (d, *J* = 8.4 Hz, 2H), 7.63 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.25, 29.41, 34.19, 40.77, 54.24, 54.84, 79.27, 113.55, 116.99, 129.35, 130.93, 133.20, 134.73, 156.15, 158.36, 172.74.

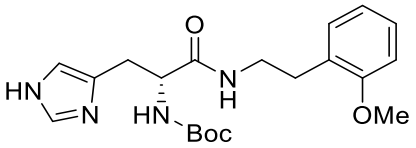


tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-((4-methoxyphenethyl)amino)-1-oxopropan-2-yl)carbamate: Following general synthetic procedure A the title compound was synthesized as a white solid (68%). R_f = 0.43 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 2.69 (t, *J* = 7.22 Hz, 2H), 2.83 (dd, *J* = 14.39, 8.93 Hz, 1H), 3.00 (dd, *J* = 14.81, 5.34 Hz, 1H), 3.35-3.44 (m, 2H), 3.77 (s, 3H), 4.25 (t, *J* = 6.59 Hz, 1H), 6.85 (d, *J* = 8.42 Hz, 3H), 7.11 (d, *J* = 8.37 Hz, 2H), 7.62 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 27.28, 29.45, 34.21, 40.78, 54.26, 54.86, 79.27, 113.56, 117.02, 129.36, 130.93, 133.20, 133.21, 134.76, 156.15, 158.35, 172.75.

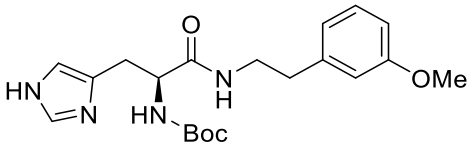


tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((2-methoxyphenethyl)amino)-1-oxopropan-2-yl)carbamate: Following general synthetic procedure A the title compound was synthesized as a white solid (70%). R_f = 0.38 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 2.80 (dt,

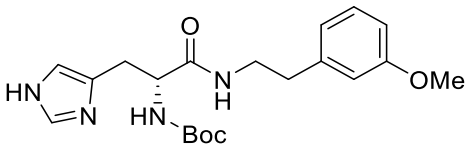
J = 23.29, 7.24 Hz, 3H), 3.00 (dd, *J* = 14.72, 5.21 Hz, 1H), 3.35 - 3.46 (m, 2H), 3.84 (s, 3H), 4.25 (t, *J* = 6.73 Hz, 1H), 6.83 - 6.88 (m, 2H), 6.93 (d, *J* = 8.11 Hz, 1H), 7.10 (d, *J* = 7.16 Hz, 1H), 7.20 (td, *J* = 7.82, 1.39 Hz, 1H), 7.60 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 25.73, 27.96, 28.16, 37.79, 52.87, 53.33, 77.72, 108.56, 115.45, 118.63, 125.46, 125.96, 128.56, 133.19, 156.05, 171.16.



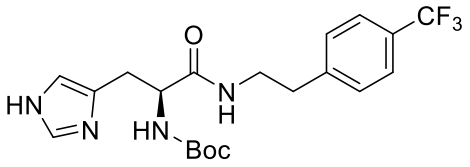
tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-((2-methoxyphenethyl)amino)-1-oxopropan-2-yl)carbamate: Following general synthetic procedure A the title compound was synthesized as a white solid (72%). R_f = 0.38 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 2.76-2.85 (m, 3H), 3.00 (dd, *J* = 14.70, 5.15 Hz, 1H), 3.35 - 3.44 (m, 2H), 3.85 (s, 3H), 4.25 (t, *J* = 6.59 Hz, 1H), 6.86 (dd, *J* = 13.33, 5.19 Hz, 2H), 6.93 (d, *J* = 8.14 Hz, 1H), 7.10 (d, *J* = 7.19 Hz, 1H), 7.20 (td, *J* = 7.82, 1.46 Hz, 1H), 7.62 (s, 1H).^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 27.27, 29.47, 29.70, 39.35, 54.42, 54.86, 79.28, 110.12, 117.03, 120.20, 127.01, 127.53, 130.12, 133.18, 134.75, 156.15, 157.61, 172.72.



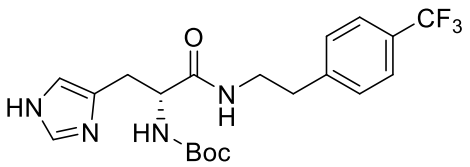
tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((3-methoxyphenethyl)amino)-1-oxopropan-2-yl)carbamate: Following general synthetic procedure A the title compound was synthesized as a white solid (71%). R_f = 0.38 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 2.73 (t, *J* = 7.25, 2H), 2.83 (dd, *J* = 14.64, 8.67 Hz, 1H), 3.00 (dd, *J* = 14.74, 5.27 Hz, 1H), 3.36 - 3.47 (m, 2H), 3.78 (s, 3H), 4.26 (t, *J* = 6.77 Hz, 1H), 6.75 - 6.79 (m, 3H), 6.85 (s, 1H), 7.19 (t, *J* = 7.89 Hz, 1H), 7.63 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 25.71, 27.91, 33.59, 38.95, 52.65, 53.32, 77.72, 109.97, 112.42, 115.45, 119.16, 127.54, 131.80, 133.21, 138.98, 154.60, 158.36, 171.24.



tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-((3-methoxyphenethyl)amino)-1-oxopropan-2-yl)carbamate: Following general synthetic procedure A the title compound was synthesized as a white solid (69%). R_f = 0.35 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 2.72 (q, *J* = 6.78, 2H), 2.83 (dd, *J* = 14.65, 8.68 Hz, 1H), 2.99 (dd, *J* = 14.67, 5.39 Hz, 1H), 3.36 - 3.48 (m, 2H), 3.79 (s, 3H), 4.25 (t, *J* = 6.68 Hz, 1H), 6.76 - 6.79 (m, 3H), 6.84 (s, 1H), 7.19 (t, *J* = 7.91 Hz, 1H), 7.60 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 27.28, 29.51, 35.15, 40.49, 54.21, 54.89, 79.26, 111.52, 113.97, 117.01, 120.72, 129.10, 133.24, 134.78, 140.52, 156.14, 159.90, 172.80.

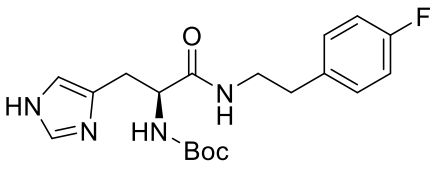


***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-(trifluoromethyl)phenethyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (84%). R_f = 0.36 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.41 (s, 9H), 2.80-2.87 (m, 3H), 2.97-3.02 (m, 1H), 3.38-3.50 (m, 2H), 4.24 (t, *J* = 6.8 Hz, 1H), 6.85 (s, 1H), 7.40 (d, *J* = 8 Hz, 2H), 7.58 (d, *J* = 8Hz, 3H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.24, 29.45, 34.81, 40.07, 54.90, 79.25, 123.10, 124.90, 124.94, 129.16, 134.79, 143.77, 156.17, 172.98.

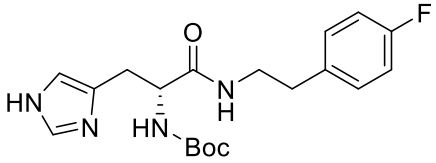


***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-(trifluoromethyl)phenethyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white

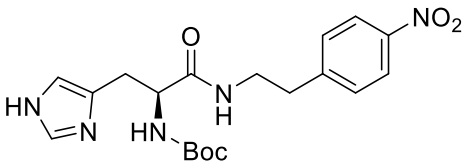
solid (70%). R_f = 0.49 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.41 (s, 9H), 2.80-2.85 (m, 2H), 2.90- 3.05 (m, 2H), 3.40-3.49 (m, 2H), 4.25 (t, *J* = 6.6 Hz, 1H), 6.86 (s, 1H), 7.40 (d, *J* = 8Hz, 2H), 7.58-7.64 (m, 3H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.23, 29.36, 34.80, 40.07, 54.84, 79.27, 116.80, 124.90, 124.93, 125.80, 129.16, 134.74, 143.77, 156.25, 172.97.



***tert-butyl (S)-(1-((4-fluorophenethyl)amino)-3-(1H-imidazol-4-yl)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (83%). R_f = 0.51 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.41 (s, 9H), 2.72-2.76 (m, 2H), 2.79-2.85 (m, 1H), 2.97-3.02 (m, 1H), 3.37-3.45 (m, 2H), 4.24 (t, *J* = 6.6 Hz, 1H), 6.84 (s, 1H), 7.00 (t, *J* = 8.6 Hz, 2H), 7.20 (t, *J* = 5.6 Hz, 2H), 7.60 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.24, 29.50, 34.21, 40.54, 54.87, 79.24, 114.55, 114.76, 130.05, 130.13, 134.77, 134.94, 160.42, 162.83, 172.87.

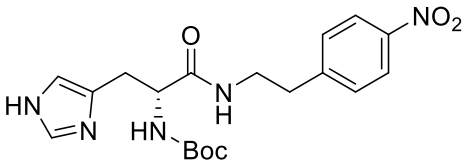


***tert-butyl (R)-(1-((4-fluorophenethyl)amino)-3-(1H-imidazol-4-yl)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (73%). R_f = 0.49 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.40 (s, 9H), 2.74 (t, *J* = 6.8 Hz, 2H), 2.82-3.37 (m, 4H), 4.27 (s, 1H), 6.90 (s, 1H), 6.99 (t, *J* = 8.4 Hz, 2H), 7.20 (t, *J* = 6.4 Hz, 2H), 7.73 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.21, 28.74, 34.18, 40.55, 54.43, 79.39, 114.6, 116.92, 130.1, 132.31, 134.63, 156.14, 161.6, 172.34.

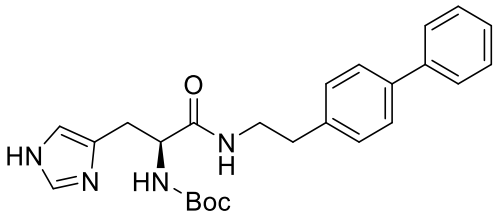


***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((4-nitrophenethyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (63%). R_f

= 0.28 (MeOH : DCM = 1 : 9).[1]H NMR (400 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 2.82 (dd, *J* = 14.74, 8.60 Hz, 1H), 2.91 (t, *J* = 6.98 Hz, 2H), 2.99 (dd, *J* = 14.76, 5.55 Hz, 1H), 3.47 (t, *J* = 6.92 Hz, 2H), 4.24 (t, *J* = 6.95 Hz, 1H), 6.86 (s, 1H), 7.45 (d, *J* = 8.48 Hz, 2H), 7.63 (s, 1H), 8.17 (d, *J* = 8.60 Hz, 2H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 27.26, 29.38, 34.83, 39.79, 54.84, 79.27, 116.84, 123.12, 129.66, 133.30, 134.75, 146.67, 147.19, 156.15, 172.96.

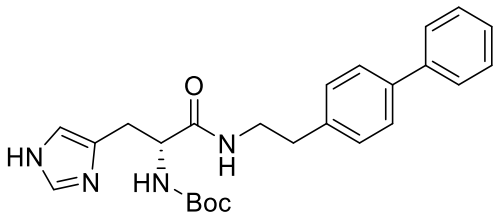


***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-((4-nitrophenethyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (64%). R_f = 0.28 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 2.82 (dd, *J* = 14.68, 8.69 Hz, 1H), 2.91 (t, *J* = 6.96 Hz, 2H), 2.99 (dd, *J* = 14.74, 5.34 Hz, 1H), 3.47 (t, *J* = 6.92 Hz, 2H), 4.24 (t, *J* = 6.93 Hz, 1H), 6.86 (s, 1H), 7.45 (d, *J* = 8.48 Hz, 2H), 7.62 (s, 1H), 8.17 (d, *J* = 8.62 Hz, 2H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 27.26, 29.41, 34.83, 39.80, 54.86, 79.27, 116.85, 123.13, 129.66, 133.33, 134.77, 146.67, 147.20, 156.15, 172.98.

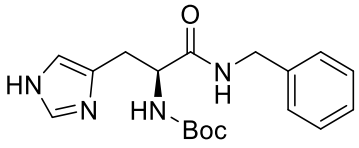


***tert-butyl (S)-(1-((2-([1,1'-biphenyl]-4-yl)ethyl)amino)-3-(1H-imidazol-4-yl)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (67%). R_f = 0.82 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.40 (s, 9H), 2.79-2.87 (m, 3H), 3.01 (dd, *J* = 14.52, 5.37 Hz, 1H), 3.38-3.52 (m, 2H), 4.27 (t, *J* = 6.82 Hz, 1H), 6.85 (s, 1H), 7.28-7.34 (m, 3H), 7.42 (t, *J* = 7.63 Hz, 2H), 7.54-7.60 (m, 5H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C :

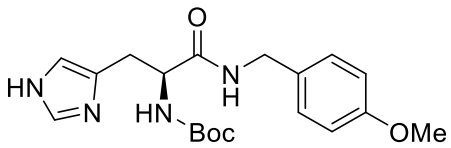
27.27, 29.50, 34.70, 40.51, 54.89, 79.26, 126.42, 126.69, 126.76, 128.41, 128.95, 134.77, 138.16, 139.18, 140.84, 172.85.



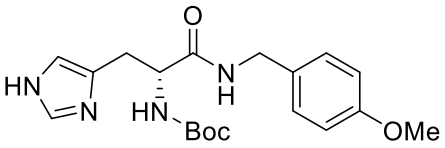
***tert-butyl (R)-(1-((2-([1,1'-biphenyl]-4-yl)ethyl)amino)-3-(1H-imidazol-4-yl)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (67%). R_f = 0.82 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 2.79-2.87 (m, 3H), 3.01 (dd, *J* = 14.61, 5.32 Hz, 1H), 3.40-3.48 (m, 2H), 4.27 (t, *J* = 6.82 Hz, 1H), 6.85 (s, 1H), 7.28-7.34 (m, 3H), 7.42 (t, *J* = 7.67 Hz, 2H), 7.55-7.60 (m, 5H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C : 27.27, 29.50, 34.70, 40.51, 54.90, 79.26, 126.42 , 126.69, 126.76, 128.41, 128.95, 134.76, 138.16, 139.18, 140.85, 156.16, 172.84.



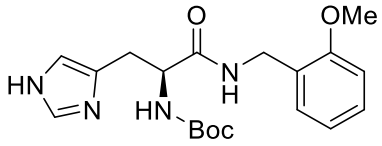
tert-butyl (S)-(1-(benzylamino)-3-(1H-imidazol-4-yl)-1-oxopropan-2-yl)carbamate: Following general synthetic procedure A the title compound was synthesized as a white solid (80%). R_f = 0.42 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.42 (s, 9H), 2.88-2.94 (m, 1H), 3.05-3.10 (m, 1H), 4.31-4.43 (m, 3H), 6.87 (s, 1H), 7.20-7.32 (m, 5H), 7.71 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.25, 29.14, 42.61, 54.88, 79.34, 116.95, 126.75, 127.01, 128.07, 132.96, 134.67, 138.35, 156.20, 172.71.



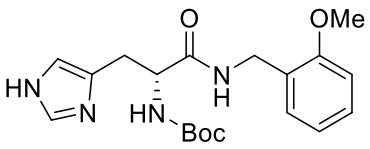
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((4-methoxybenzyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (70%). R_f = 0.50 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.42 (m, 9H), 2.89 (dd, *J* = 14.44, 8.39 Hz, 1H), 3.05 (dd, *J* = 14.55, 5.74 Hz, 1H), 3.78 (s, 3H), 4.24-4.35 (m, 3H), 6.85 (d, *J* = 8.77 Hz, 3H), 7.13 (d, *J* = 8.19 Hz, 2H), 7.62 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 27.26, 29.36, 42.12, 54.29, 54.96, 79.29, 113.45, 116.98, 128.36, 130.31, 133.19, 134.76, 156.18, 158.97, 172.66.



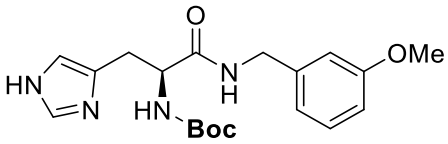
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-(trifluoromethyl)phenethyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (77%). R_f = 0.31 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.41 (s, 9H), 2.85-3.06 (m, 2H), 3.78 (s, 3H), 4.29 (m, 3H), 6.83-6.86 (m, 3H), 7.13 (d, *J* = 8 Hz, 2H), 7.59 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.24, 29.38, 42.10, 54.27, 54.98, 79.27, 113.44, 128.36, 130.30, 134.78, 156.27, 158.98, 172.711.



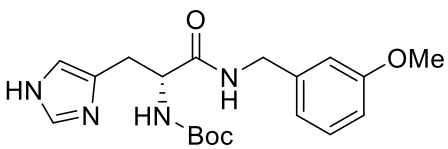
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((2-methoxybenzyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (75%). R_f = 0.46 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 2.87-2.91 (m, 1H), 3.01-3.07 (m, 1H), 3.85 (s, 3H), 4.32-4.42 (m, 3H), 6.81 (s, 1H), 6.88 (t, *J* = 7.42 Hz, 1H), 6.95 (d, *J* = 8.17 Hz, 1H), 7.11 (d, *J* = 7.27 Hz, 1H), 7.25 (t, *J* = 7.63 Hz, 1H), 7.60 (s, 1H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C : 27.26, 29.34, 38.15, 54.43, 54.94, 79.29, 109.92, 117.00, 120.00, 125.81, 128.03, 128.19, 134.74, 156.22, 157.29, 172.75.



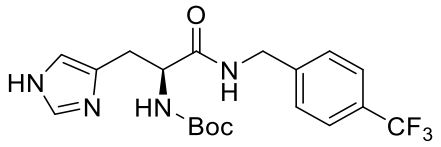
***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-((2-methoxybenzyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (65%). R_f = 0.43 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H :1.42 (m, 9H), 2.88 (dd, *J* = 14.64, 8.44 Hz, 1H), 3.05 (dd, *J* = 14.67, 5.51 Hz, 1H), 3.85 (s, 3H), 4.31-4.42 (m, 3H), 6.81 (s, 1H), 6.88 (t, *J* = 7.40 Hz, 1H), 6.95 (d, *J* = 8.18 Hz, 1H), 7.12 (d, *J* = 7.30 Hz, 1H), 7.25 (t, *J* = 7.16 Hz, 1H), 7.58 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 27.29, 29.38, 38.16, 54.44, 54.96, 79.29, 109.91, 120.00, 125.81, 128.01, 128.19, 134.76, 156.21, 157.27, 172.77.



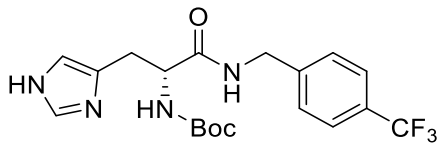
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((3-methoxybenzyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (65%). R_f = 0.47 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H : 1.42(s, 9H), 2.88-2.93 (m, 1H), 3.01-3.08 (m, 1H), 3.79 (s, 3H), 4.34 (q, *J* = 13.41 Hz, 3H), 6.78-6.85 (m, 4H), 7.21 (t, *J* = 7.77 Hz, 1H), 7.61 (s, 1H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C : 27.27, 29.36, 42.55, 54.25, 55.01, 79.28, 112.30, 112.50, 119.17, 129.06, 134.78, 139.89, 159.93, 172.88.



***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-((3-methoxybenzyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (61%). R_f = 0.42 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 2.90 (dd, *J* = 14.61, 8.45 Hz, 1H), 3.06 (dd, *J* = 14.70, 5.65 Hz, 1H), 3.79 (m, 3H), 4.30- 4.39 (m, 3H), 6.81 (dd, *J* = 18.07, 7.73 Hz, 4H), 7.21 (t, *J* = 7.79 Hz, 1H), 7.61 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 27.26, 29.34, 42.55, 54.25, 55.00, 79.29, 112.30, 112.50, 116.91, 119.19, 129.06, 133.29, 134.77, 139.90, 156.20, 159.93, 172.90.

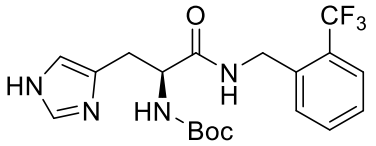


***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-(trifluoromethyl)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (74%). R_f = 0.39 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.43 (s, 9H), 2.89-2.95 (m,1H), 3.01-3.09 (m, 1H), 4.33-4.50 (m, 3H), 6.87 (s, 1H), 7.38 (d, *J* = 7.6 Hz, 2H), 7.58-7.62 (m, 3H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.24, 29.24, 42.06, 55.08, 79.32, 116.96, 123.22, 124.87, 125.64, 127.19, 127.38, 129.04, 134.85, 143.15, 156.21,173.16.

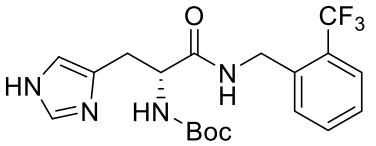


***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-(trifluoromethyl)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (69%). R_f = 0.31 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.41 (s, 9H), 2.93-3.08 (m, 2H), 4.36-4.44 (m, 3H), 6.93 (s, 1H), 7.39 (s, 2H), 7.58 (s, 2H), 7.80 (s, 1H); ^{13}C NMR (100 MHz,

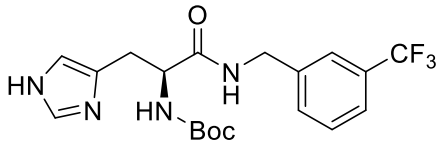
MeOH-d_4) δ_C: 27.23, 28.86, 42.09, 54.87, 79.41, 116.89, 122.98, 124.88, 125.67, 127.41, 132.86, 134.65, 143.12, 156.23, 172.91.



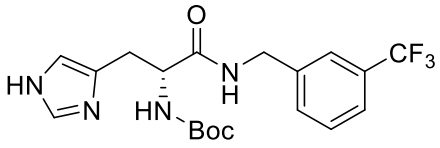
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((2-(trifluoromethyl)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (64%). R_f = 0.45 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 2.92 (dd, *J* = 14.61, 8.18 Hz, 1H), 3.06 (dd, *J* = 14.65, 5.96 Hz, 1H), 4.36 (t, *J* = 6.90 Hz, 1H), 4.56 (q, *J* = 12.57 Hz, 2H), 6.86 (s, 1H), 7.35-7.43 (m, 2H), 7.54 (t, *J* = 7.55 Hz, 1H), 7.63 - 7.67 (m, 2H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 25.72, 27.62, 37.60, 53.54, 77.81, 115.34, 121.66, 123.83, 123.89, 124.37, 125.46, 126.86, 130.49, 133.32, 135.07, 154.74, 171.69.



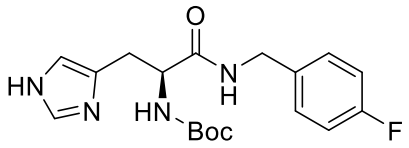
***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-((2-(trifluoromethyl)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (72%). R_f = 0.45 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H : 1.43 (s, 9H), 2.95 (dd, *J* = 14.60, 8.19 Hz, 1H), 3.09 (dd, *J* = 14.62, 5.88 Hz, 1H), 4.40 (t, *J* = 6.91 Hz, 1H), 4.58 (q, *J* = 15.54 Hz, 2H), 6.88 (s, 1H), 7.37-7.43 (m, 2H), 7.54 (t, *J* = 7.50 Hz, 1H), 7.63 (s, 1H), 7.67 (d, *J* = 7.74 Hz, 1H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C : 27.30, 29.24, 39.14, 55.13, 79.36, 116.91, 121.29, 123.46, 125.33, 125.37, 125.43, 125.63, 126.99, 127.06, 127.31, 127.81, 128.39, 132.03, 133.40, 134.89, 136.61, 156.27, 173.27.



***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((3-(trifluoromethyl)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (82%). R_f = 0.33 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H: 1.42 (s, 9H), 2.88-2.93 (m, 1H), 3.04-3.08 (m, 1H), 4.32-4.46 (m, 3H), 6.84 (s, 1H), 7.47-7.60 (m, 5H); ^{13}C NMR (125 MHz, MeOH-d_4) δ_C: 27.2, 29.6, 42.1, 55.0, 79.3, 116.8, 123.1, 123.4, 123.7, 125.3, 130.2, 130.4, 130.7, 134.8, 140.0, 156.2, 173.1.

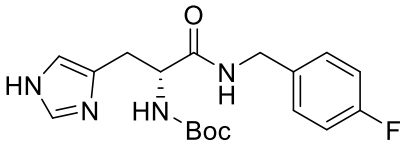


***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-((3-(trifluoromethyl)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (74%). R_f = 0.35 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, DMSO-d_6) δ_H: 1.42 (s, 9H), 2.88-2.93 (m, 1H), 3.04-3.08 (m, 1H), 4.32-4.47 (m, 3H), 6.84 (s, 1H), 7.47-7.61 (m, 5H); ^{13}C NMR (100 MHz, DMSO-d_6) δ_C: 27.2, 29.2, 42.0, 55.0, 79.3, 99.9, 123.1, 123.4, 123.7, 125.3, 128.3, 130.2, 130.4, 130.7, 134.8, 139.9, 156.2, 173.1.

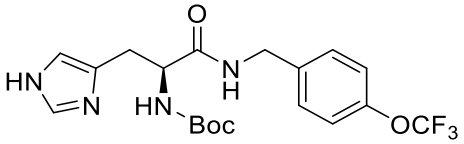


***tert-butyl (S)-(1-((4-fluorobenzyl)amino)-3-(1H-imidazol-4-yl)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (86%). R_f = 0.34 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 2.87-2.92 (m, 1H), 3.02-

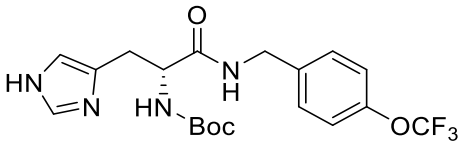
3.07 (m, 1H), 4.34 (q, *J* = 14.3 Hz, 3H), 6.84 (s, 1H), 7.02 (t, *J* = 8.8 Hz, 2H), 7.21 (t, *J* = 6.6 Hz), 7.61 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.24, 29.29, 41.84, 55.01, 79.30, 114.62, 128.93, 134.80, 163.34, 172.88.



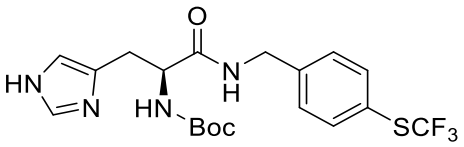
***tert-butyl (R)-(1-((4-fluorobenzyl)amino)-3-(1H-imidazol-4-yl)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (70 %). R_f = 0.33 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 2.94-3.00 (m, 1H), 3.08-3.22 (m, 1H), 4.27-4.40 (m, 3H), 7.04 (t, *J* = 8.6 Hz, 2H), 7.14 (s, 1H), 7.26-7.29 (m, 2H), 8.43 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.18, 27.71, 35.60, 41.96, 54.05, 79.66, 114.7, 116.95, 128.98, 129.08, 131.08, 133.89, 134.43, 163.39, 174.42.



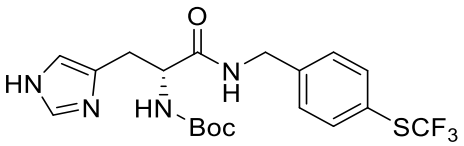
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-(trifluoromethoxy)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (89%). R_f = 0.38 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 2.87-2.93 (m, 1H), 3.03-3.08 (m, 1H), 4.31-4.43 (m, 3H), 6.85 (s, 1H), 7.20 (d, *J* = 8.4 Hz, 2H), 7.29 (d, *J* = 8.4 Hz, 2H), 7.60 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.25, 29.29, 41.80, 55.08, 79.31, 119.26, 120.61, 121.79, 128.61, 134.85, 137.83, 148.11, 156.23, 173.05.



***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-(trifluoromethoxy)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (91%). R_f = 0.29 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 2.87-2.93 (m, 1H), 3.01-3.08 (m, 1H), 4.31-4.43 (m, 3H), 6.85 (s, 1H), 7.20 (d, *J* = 8 Hz, 2H), 7.29 (d, *J*= 8.4 Hz, 2H), 7.60 (s, 1H); ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.23, 29.26, 41.79, 55.07, 79.30, 119.25, 120.60, 121.79, 128.61, 134.83, 137.83, 148.10, 156.22, 173.04.

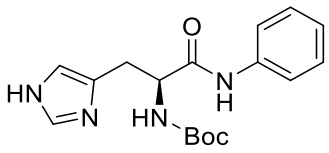


***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-((trifluoromethyl)thio)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (66%). R_f = 0.53 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.43 (s, 9H), 2.92 (dd, *J* = 14.60, 8.27 Hz, 1H), 3.07 (dd, *J* = 14.68, 5.89 Hz, 1H), 4.35 (dd, *J* = 13.14, 5.62 Hz, 1H), 4.42- 4.48 (m, 2H), 6.87 (s, 1H), 7.33 (d, *J* = 7.92 Hz, 2H), 7.62 (d, *J* = 6.18 Hz, 3H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.26, 29.21, 42.01, 55.08, 79.34, 116.88, 122.17, 128.16, 128.25, 131.30, 134.82, 136.16, 142.35, 156.24, 173.12.

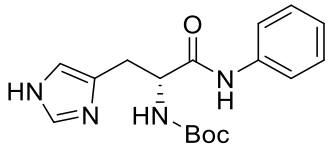


***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-((trifluoromethyl)thio)benzyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (64%). R_f = 0.53 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H: 1.43 (s, 9H), 2.92 (dd, *J* = 14.66, 8.20 Hz, 1H), 3.07 (dd, *J* = 14.64, 5.93 Hz, 1H), 4.35 (t, *J* = 6.98 Hz, 1H), 4.43 (q, *J* = 13.72 Hz, 2H), 6.88 (s, 1H), 7.33 (d, *J* = 7.92 Hz, 2H), 7.62-7.66 (m, 3H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C:

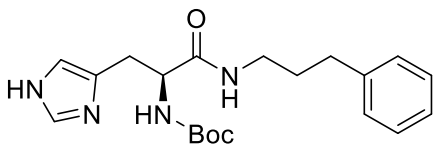
27.28, 29.19, 42.02, 55.07, 79.35, 116.91, 122.17, 128.17, 128.56, 131.00, 133.28, 134.82, 136.16, 142.35, 156.24, 173.10.



***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-(phenylamino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (64%). R_f = 0.30 (MeOH : DCM = 1 : 9).^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.43 (s, 9H), 2.96 (dd, *J* = 14.38, 8.28 Hz, 1H), 3.10 (dd, *J* = 14.50, 5.68 Hz, 1H), 4.43 (t, *J* = 6.62 Hz, 1H), 6.90 (s, 1H), 7.11 (t, *J* = 7.41 Hz, 1H), 7.31 (t, *J* = 7.89 Hz, 2H), 7.52 (d, *J* = 7.87 Hz, 2H), 7.63 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 25.73, 28.13, 53.94, 77.80, 115.27, 118.63, 122.51, 126.85, 133.29, 136.47, 154.76, 169.77.

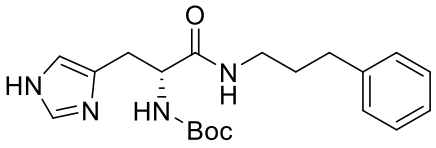


***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-(phenylamino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (68%). R_f = 0.30 (MeOH : DCM = 1 : 9).^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.44 (s, 9H), 2.96 (dd, *J* = 14.59, 8.28 Hz, 1H), 3.10 (dd, *J* = 14.73, 5.59 Hz, 1H), 4.43 (t, *J* = 6.75 Hz, 1H), 6.90 (s, 1H), 7.11 (t, *J* = 7.42 Hz, 1H), 7.31 (t, *J* = 7.92 Hz, 2H), 7.53 (d, *J* = 7.73 Hz, 2H), 7.63 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.27, 29.67, 55.47, 79.34, 116.81, 120.16, 124.04, 128.39, 133.31, 134.82, 138.00, 156.30, 171.30.



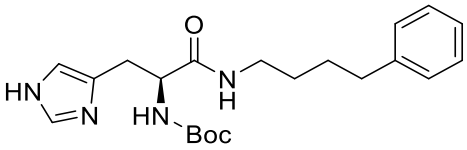
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((3-phenylpropyl)amino)propan-2-yl)carbamate*:**

Following general synthetic procedure A the title compound was synthesized as a white solid (55%). R_f = 0.53 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 1.73-1.81 (m, 2H), 2.59 (t, *J* = 7.76 Hz, 2H), 2.88 (dd, *J* = 14.68, 8.34 Hz, 1H), 3.03 (dd, *J* = 14.63, 5.65 Hz, 1H), 3.14-3.26 (m, 2H), 4.28 (t, *J* = 6.85 Hz, 1H), 6.88 (s, 1H), 7.14-7.19 (m, 3H), 7.26 (t, *J* = 7.44 Hz, 2H), 7.61 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.29, 29.46, 30.81, 32.61, 38.57, 54.98, 79.27, 116.98, 125.49, 127.99, 128.04, 133.28, 134.78, 141.60, 156.16, 172.85.



***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-((3-phenylpropyl)amino)propan-2-yl)carbamate*:**

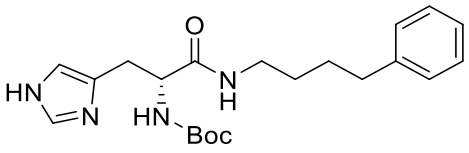
Following general synthetic procedure A the title compound was synthesized as a white solid (54%). R_f = 0.53 (MeOH : DCM = 1 : 9).^{1}H NMR (500 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 1.74-1.79 (m, 2H), 2.58 (t, *J* = 7.68 Hz, 2H), 2.89 (dd, *J* = 14.59, 8.36 Hz, 1H), 3.04 (dd, *J* = 14.66, 5.68 Hz, 1H), 3.14-3.24 (m, 2H), 4.30 (t, *J* = 6.83 Hz, 1H), 6.88 (s, 1H), 7.14-7.18 (m, 3H), 7.25 (t, *J* = 7.52 Hz, 2H), 7.61 (s, 1H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C: 27.36, 29.52, 30.81, 32.63, 38.60, 55.00, 79.27, 117.05, 125.52, 128.02, 128.07, 133.29, 134.82, 141.59, 156.16, 172.85.



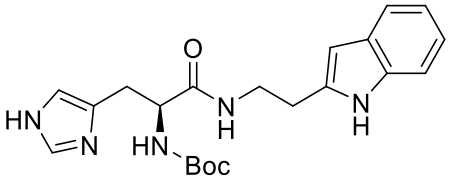
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-phenylbutyl)amino)propan-2-yl)carbamate*:**

Following general synthetic procedure A the title compound was synthesized as a white solid (51%). R_f = 0.42 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 1.46-1.52 (m, 2H), 1.57-1.63 (m, 2H), 2.61 (t, *J* = 7.50 Hz, 2H), 2.86 (dd, *J* = 14.68, 8.49 Hz, 1H), 3.02 (dd, *J* = 14.61, 5.51 Hz,

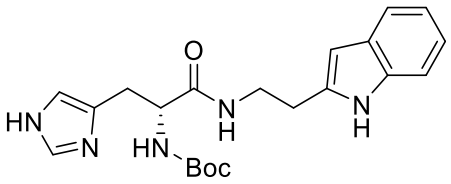
1H), 3.14-3.25 (m, 2H), 4.26 (t, *J* = 6.77 Hz, 1H), 6.86 (s, 1H), 7.14-7.19 (m, 3H), 7.25 (t, *J* = 7.53 Hz, 2H), 7.59 (s, 1H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C: 27.29, 28.35, 28.54, 29.49, 35.07, 38.74, 54.95, 79.25, 117.00, 125.35, 127.91, 127.98, 128.06, 133.25, 134.76, 142.14, 156.14, 172.77.



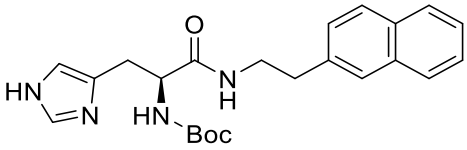
***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-((4-phenylbutyl)amino)propan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (50%). R_f = 0.42 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H : 1.41 (s, 9H), 1.46-1.52 (m, 2H), 1.57-1.63 (m, 2H), 2.61 (t, *J* = 7.50 Hz, 2H), 2.86 (q, *J* = 7.13 Hz, 1H), 3.00-3.04 (m, 1H), 3.14-3.25 (m, 2H), 4.27 (t, *J* = 6.81 Hz, 1H), 6.87 (s, 1H), 7.13-7.19 (m, 3H), 7.24-7.27 (m, 2H), 7.61 (s, 1H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C: 27.30, 28.36, 28.54, 29.44, 35.07, 38.75, 54.93, 79.26, 117.00, 125.35, 127.91, 128.07, 133.18, 134.74, 142.13, 156.14, 172.73.



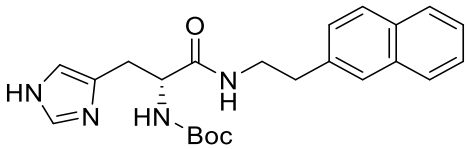
***tert-butyl (S)-(1-((2-(1H-indol-2-yl)ethyl)amino)-3-(1H-imidazol-4-yl)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a yellowish solid (75%). R_f = 0.35 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.40 (s, 9H), 2.81-2.87 (m, 1H), 3.02 (dd, *J* = 14.86, 5.26 Hz, 1H), 3.43-3.54 (m, 2H), 4.26-4.29 (m, 1H), 6.86 (s, 1H), 7.00-7.12 (m, 3H), 7.33-7.35 (m, 1H), 7.57 (d, *J* = 7.85 Hz, 1H), 7.68 (s, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C : 24.77, 27.24, 29.26, 39.85, 54.80, 79.32, 110.85, 111.62, 117.08, 117.86, 118.24, 120.95, 122.08, 127.29, 132.93, 134.63, 136.79, 156.16, 172.61.



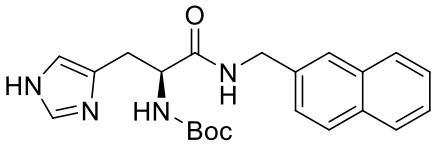
***tert-butyl (R)-(1-((2-(1H-indol-2-yl)ethyl)amino)-3-(1H-imidazol-4-yl)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a yellowish solid (73%). R_f = 0.35 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.40 (s, 9H), 2.81-2.86 (m, 1H), 2.96-3.03 (m, 1H), 3.42-3.54 (m, 2H), 4.25-4.28 (m, 1H), 6.83 (s, 1H), 7.00-7.12 (m, 3H), 7.33-7.35 (m, 1H), 7.55-7.58 (m, 2H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 23.24, 25.71, 27.91, 38.29, 53.37, 77.75, 109.31, 110.09, 116.33, 116.70, 119.42, 120.55, 125.76, 131.63, 133.21, 135.25, 154.63, 171.21.



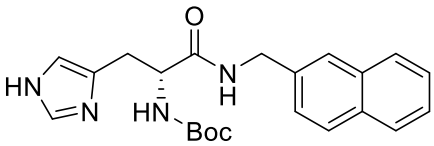
***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((2-(naphthalen-2-yl)ethyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (60%). R_f = 0.60 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H :1.37 (s, 9H), 2.81 (dd, *J* = 14.74, 8.61 Hz, 1H), 2.91-3.01 (m, 3H), 3.44-3.57 (m, 2H), 4.25 (t, *J* = 6.82 Hz, 1H), 6.82 (s, 1H), 7.36 (d, *J* = 8.50 Hz, 1H), 7.40-7.46 (m, 2H), 7.58 (s, 1H), 7.67 (s, 1H), 7.80 (td, *J* = 4.45, 2.81 Hz, 3H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 15.06, 27.23, 29.44, 35.18, 40.40, 54.86, 79.28, 116.98, 125.00, 125.58, 126.74, 126.92, 127.18, 127.71, 132.37, 133.71, 134.74, 136.50, 156.19, 172.87.



***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-((2-(naphthalen-2-yl)ethyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (61%). R_f = 0.60 (MeOH : DCM = 1 : 9).^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.38 (s, 9H), 2.81 (dd, *J* = 14.61, 8.50 Hz, 1H), 2.92-3.02 (m, 3H), 3.45-3.58 (m, 2H), 4.25 (t, *J* = 6.92 Hz, 1H), 6.83 (s, 1H), 7.37 (dd, *J* = 8.47, 1.14 Hz, 1H), 7.40-7.47 (m, 2H), 7.59 (s, 1H), 7.67 (s, 1H), 7.80 (dt, *J* = 6.67, 2.28 Hz, 3H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.22, 29.42, 35.17, 40.39, 54.87, 79.25, 116.98, 125.00, 125.58, 126.73, 126.92, 127.18, 127.71, 132.37, 133.72, 134.73, 136.51, 156.16, 172.85.

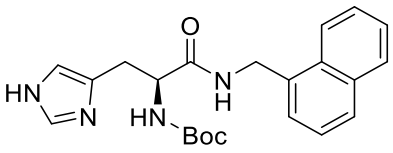


***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((naphthalen-2-ylmethyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (72%). R_f = 0.80 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 2.94 (dd, *J* = 14.60, 8.38 Hz, 1H), 3.09 (dd, *J* = 14.68, 5.66 Hz, 1H), 4.39 (t, *J* = 6.84 Hz, 1H), 4.54 (q, *J* = 16.76 Hz, 2H), 6.87 (s, 1H), 7.33 (d, *J* = 8.35 Hz, 1H), 7.43-7.48 (m, 2H), 7.60 (s, 1H), 7.70 (s, 1H), 7.79-7.83 (m, 3H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C: 27.08, 27.29, 29.38, 42.75, 55.09, 79.31, 116.96, 125.28, 125.35, 125.39, 125.73, 127.24, 127.36, 127.80, 132.76, 133.47, 134.82, 135.81, 156.26, 173.04.

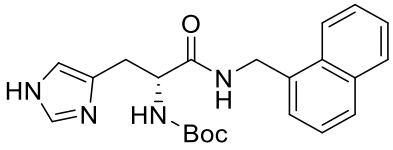


***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-((naphthalen-2-ylmethyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (70%). R_f = 0.80 (MeOH : DCM = 1 : 9). ^{1}H NMR (500 MHz, MeOH-d_4) δ_H : 1.42 (s, 9H), 2.94 (dd, *J* = 14.61, 8.29 Hz, 1H), 3.09 (dd, *J* = 14.62, 5.70 Hz, 1H), 4.39 (t, *J* = 6.86 Hz, 1H), 4.54 (q, *J* = 16.77

Hz, 2H), 6.87 (s, 1H), 7.33 (d, *J* = 8.32 Hz, 1H), 7.43-7.48 (m, 2H), 7.60 (s, 1H), 7.70 (s, 1H), 7.79-7.83 (m, 3H). ^{13}C NMR (126 MHz, MeOH-d_4) δ_C: 27.29, 29.39, 42.75, 55.09, 79.31, 117.11, 125.27, 125.35, 125.39, 125.72, 127.23, 127.35, 127.79, 132.75, 133.46, 134.82, 135.81, 156.25, 173.03.

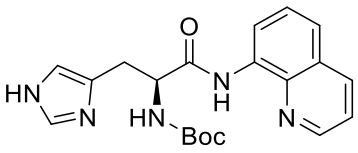


***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-((naphthalen-1-ylmethyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (76%). R_f = 0.50 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.39 (s, 9H), 2.90 (dd, *J* = 14.60, 8.36 Hz, 1H), 3.04-3.08 (m, 1H), 4.37 (t, *J* = 6.87 Hz, 1H), 4.83 (q, *J* = 21.08 Hz, 2H), 6.83 (s, 1H), 7.39 (dt, *J* = 18.35, 8.43 Hz, 2H), 7.48-7.54 (m, 2H), 7.58 (s, 1H), 7.80 (d, *J* = 8.0 Hz, 1H), 7.88 (d, *J* = 7.5 Hz, 1H), 7.99 (d, *J* = 7.76 Hz, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.27, 29.40, 40.75, 54.99, 79.28, 116.90, 122.95, 124.98, 125.29, 125.32, 125.34, 125.43, 125.98, 127.75, 128.32, 131.23, 133.23, 133.30, 133.89, 134.75, 156.19, 172.73.



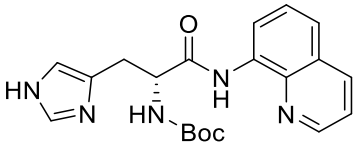
***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-((naphthalen-1-ylmethyl)amino)-1-oxopropan-2-yl)carbamate*:** Following general synthetic procedure A the title compound was synthesized as a white solid (75%). R_f = 0.50 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H: 1.39 (s, 9H), 2.90 (dd, *J* = 14.60, 8.34 Hz, 1H), 3.05(dd, *J* = 14.63, 5.61 Hz, 1H), 4.36 (t, *J* = 6.91 Hz, 1H), 4.82 (t, *J* = 16.60 Hz, 2H), 6.82 (s, 1H), 7.37-7.44 (m, 2H), 7.48-7.58 (m, 3H), 7.81 (d, *J* = 7.95 Hz, 1H), 7.89 (dd, *J* = 7.21, 2.11 Hz, 1H), 8.00 (d, *J* = 7.82 Hz, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.00, 27.25, 29.43,

40.74, 55.01, 79.28, 116.92, 122.96, 124.99, 125.33, 125.35, 125.44, 125.99, 127.75, 128.32, 131.24, 133.31, 133.90, 134.78, 156.22, 172.78.



***tert-butyl (S)-(3-(1H-imidazol-4-yl)-1-oxo-1-(quinolin-8-ylamino)propan-2-yl)carbamate*:**

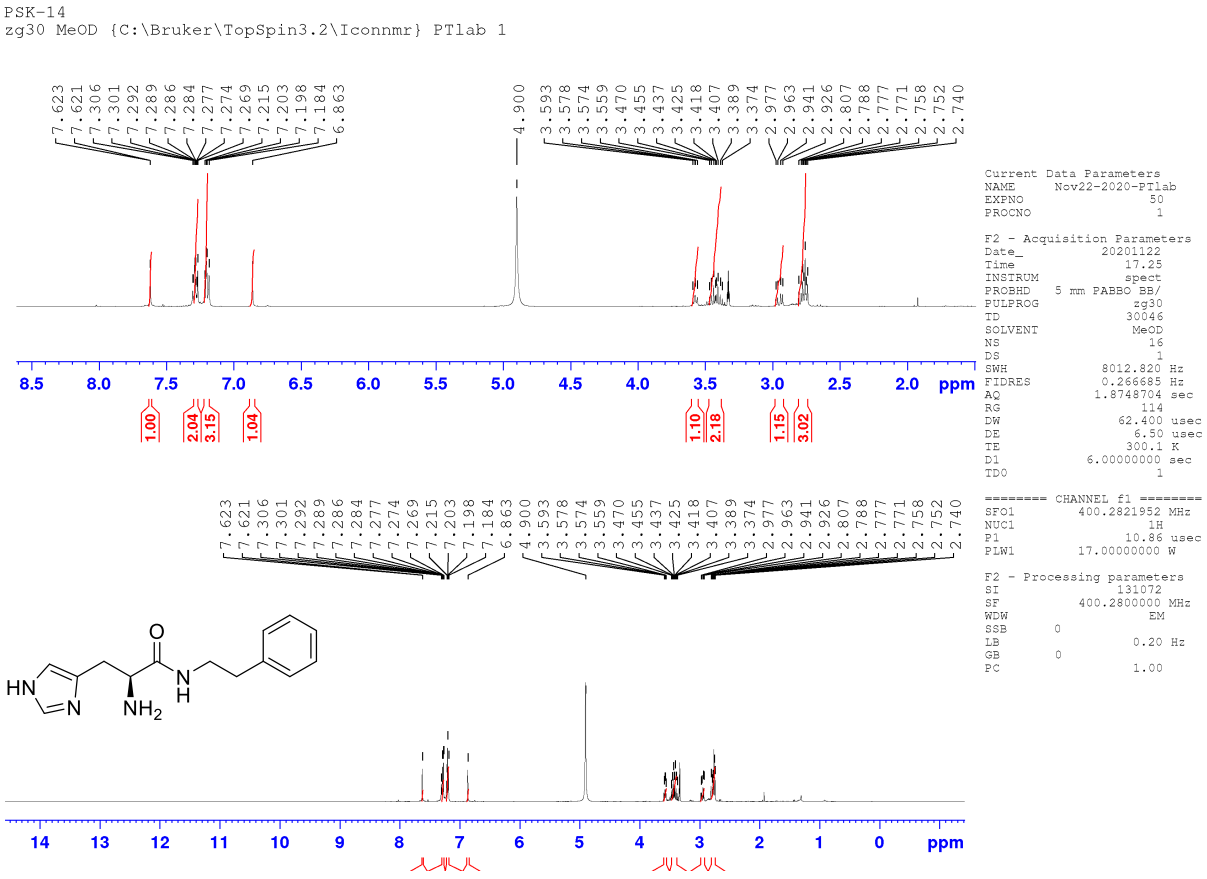
Following general synthetic procedure A the title compound was synthesized as a yellowish solid (56%). R_f = 0.53 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.48 (s, 9H), 3.07 (dd, *J* = 14.87, 9.10 Hz, 1H), 3.29 (d, *J* = 4.75 Hz, 1H), 4.59 (dd, *J* = 8.86, 5.15 Hz, 1H), 6.93 (s, 1H), 7.54-7.58 (m, 2H), 7.64 (dd, *J* = 9.25, 2.05 Hz, 2H), 8.30 (dd, *J* = 8.32, 1.60 Hz, 1H), 8.68 (dd, *J* = 7.60, 1.35 Hz, 1H), 8.84 (dd, *J*=4.24, 1.66 Hz, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.33, 28.70, 56.18, 79.67, 116.19, 121.70, 122.11, 126.59, 128.12, 133.68, 134.95, 136.15, 138.56, 148.47, 156.56, 171.00.



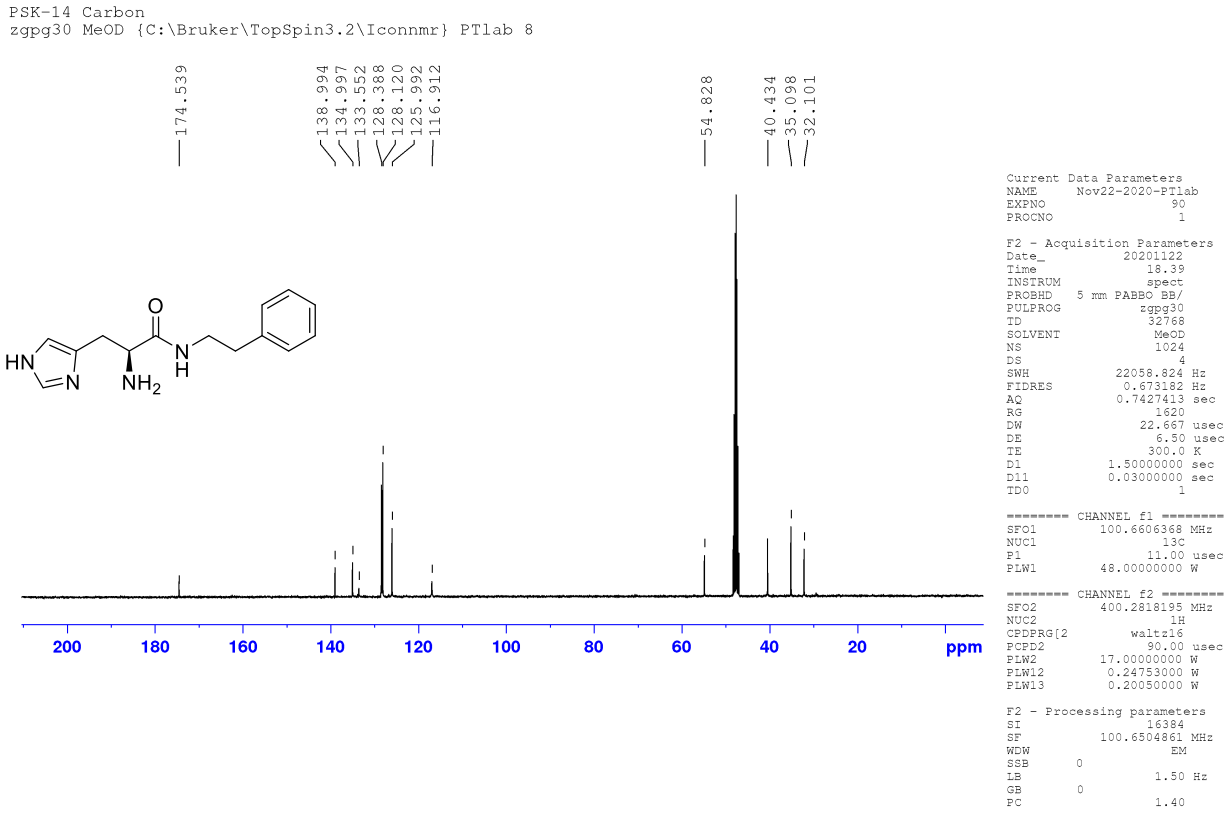
***tert-butyl (R)-(3-(1H-imidazol-4-yl)-1-oxo-1-(quinolin-8-ylamino)propan-2-yl)carbamate*:**

Following general synthetic procedure A the title compound was synthesized as a yellowish solid (54%). R_f = 0.53 (MeOH : DCM = 1 : 9). ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : ^{1}H NMR (400 MHz, MeOH-d_4) δ_H : 1.48 (s, 9H), 3.07 (dd, *J* = 14.69, 8.99 Hz, 1H), 3.28 (d, *J* = 4.83 Hz, 1H), 4.59 (dd, *J* = 8.62, 5.10 Hz, 1H), 6.93 (s, 1H), 7.54-7.58 (m, 2H), 7.63 (d, *J* = 5.26 Hz, 2H), 8.30 (dd, *J* = 8.31, 1.54 Hz, 1H), 8.68 (dd, *J* = 7.59, 1.28 Hz, 1H), 8.84 (dd, *J* = 4.22, 1.62 Hz, 1H). ^{13}C NMR (100 MHz, MeOH-d_4) δ_C: 27.35, 28.79, 56.23, 79.66, 116.18, 117.12, 121.68, 122.09, 126.57, 128.08, 133.41, 133.66, 134.98, 136.12, 138.53, 148.44, 156.54, 171.02.

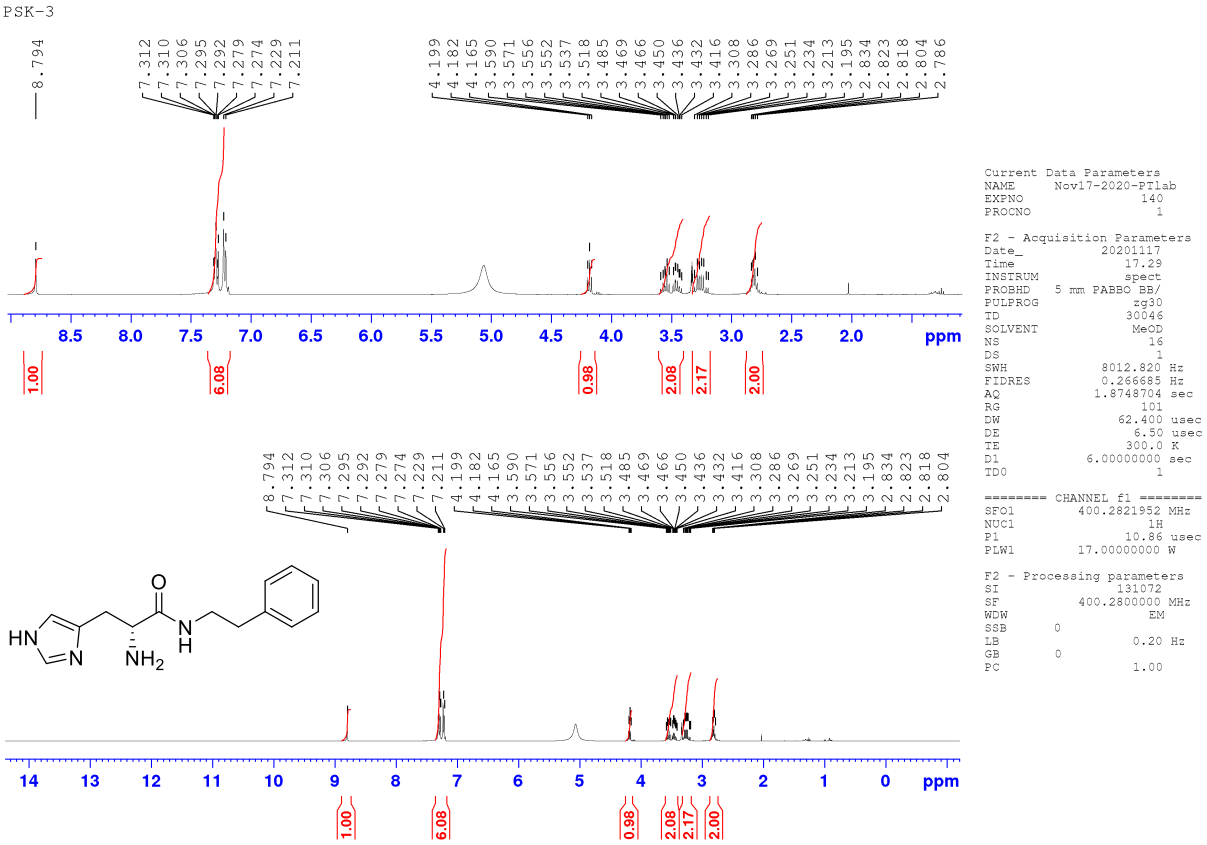
^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenethylpropanamide **(4c):**



^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenethylpropanamide **(4c):**



^{1}H NMR (400 MHz, MeOH-d$_4$), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenethylpropanamide **(4d):**

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Current Data Parameters
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EXPNO              140
PROCNO               1

F2 - Acquisition Parameters
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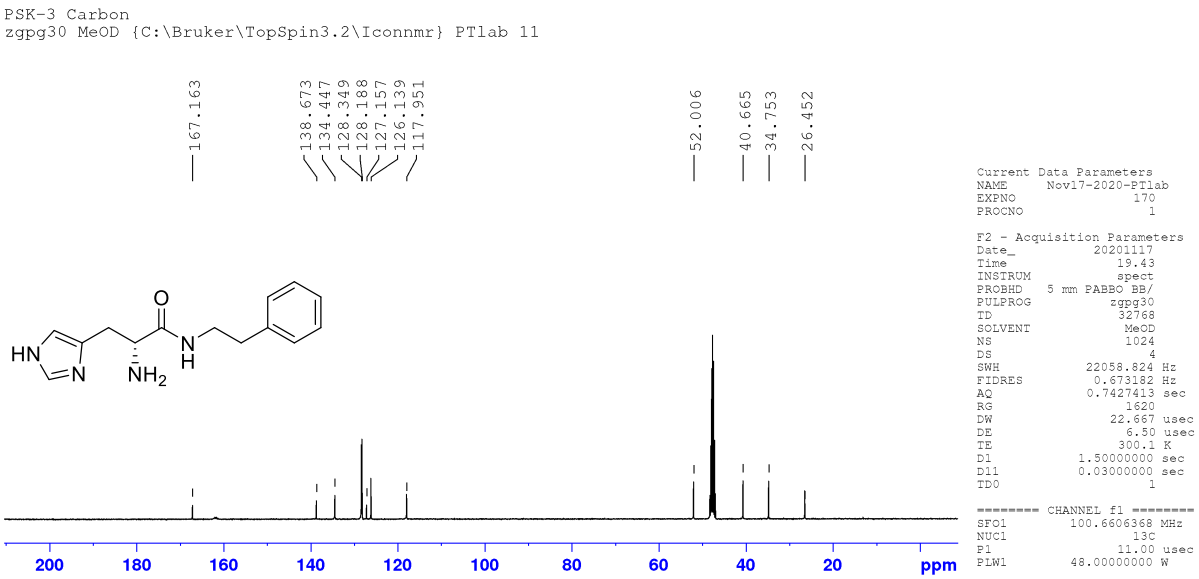
Time              17.29
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG           zg30
TD               30046
SOLVENT           MeOD
NS                  16
DS                   1
SWH           8012.820 Hz
FIDRES        0.266685 Hz
AQ            1.8748704 sec
RG                 101
DW              62.400 usec
DE               6.50 usec
TE              300.0 K
D1            6.00000000 sec
TD0                  1

======== CHANNEL f1 ========
SFO1        400.2821952 MHz
NUC1               1H
P1               10.86 usec
PLW1       17.00000000 W

F2 - Processing parameters
SI              131072
SF          400.2800000 MHz
WDW                EM
SSB       0
LB               0.20 Hz
GB        0
PC               1.00
```

^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenethylpropanamide **(4d):**

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Current Data Parameters
NAME      Nov17-2020-PTlab
EXPNO                  170
PROCNO                   1

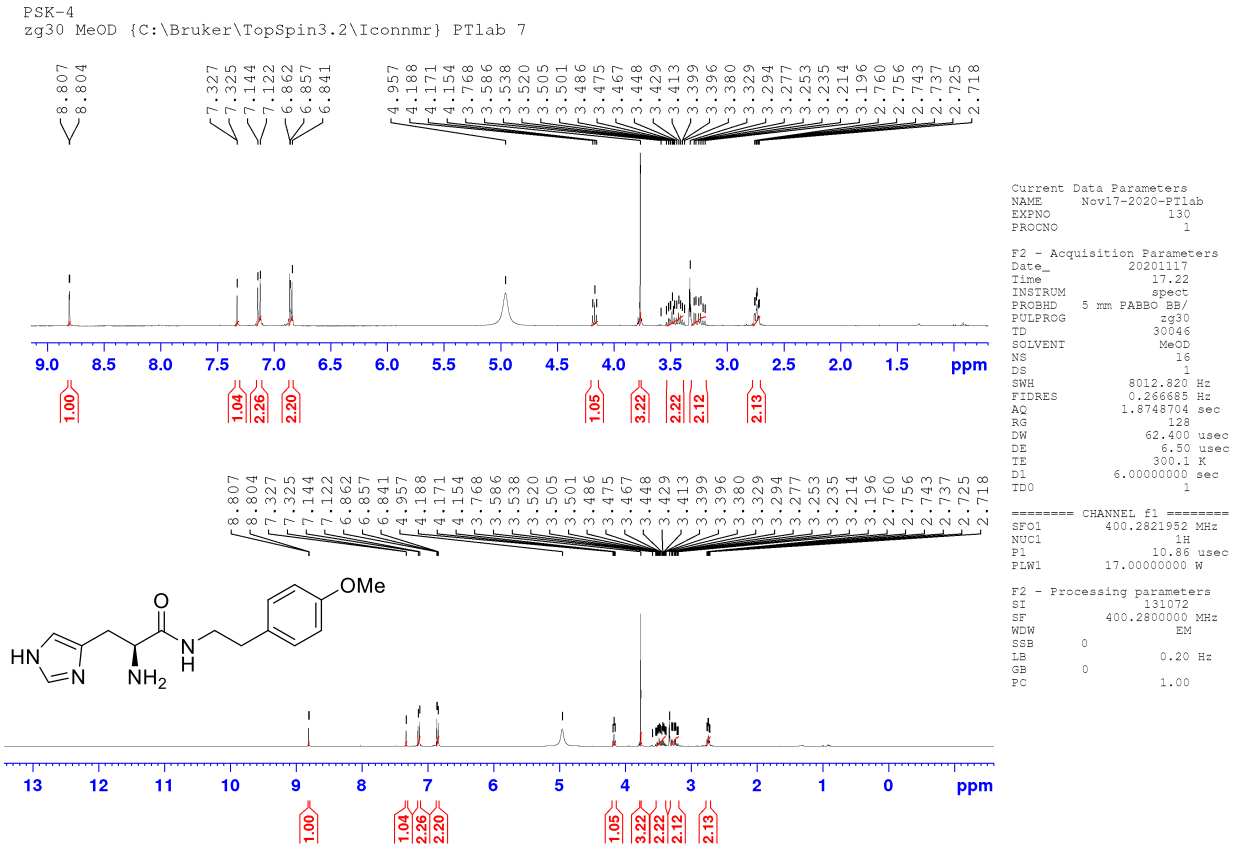
F2 - Acquisition Parameters
Date_           20201117
Time              19.43
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG         zgpg30
TD               32768
SOLVENT           MeOD
NS                1024
DS                   4
SWH           22058.824 Hz
FIDRES         0.673182 Hz
AQ            0.7427413 sec
RG                1620
DW              22.667 usec
DE               6.50 usec
TE              300.1 K
D1           1.50000000 sec
D11          0.03000000 sec
TD0                  1

======== CHANNEL f1 ========
SFO1        100.6606368 MHz
NUC1              13C
P1              11.00 usec
PLW1        48.00000000 W

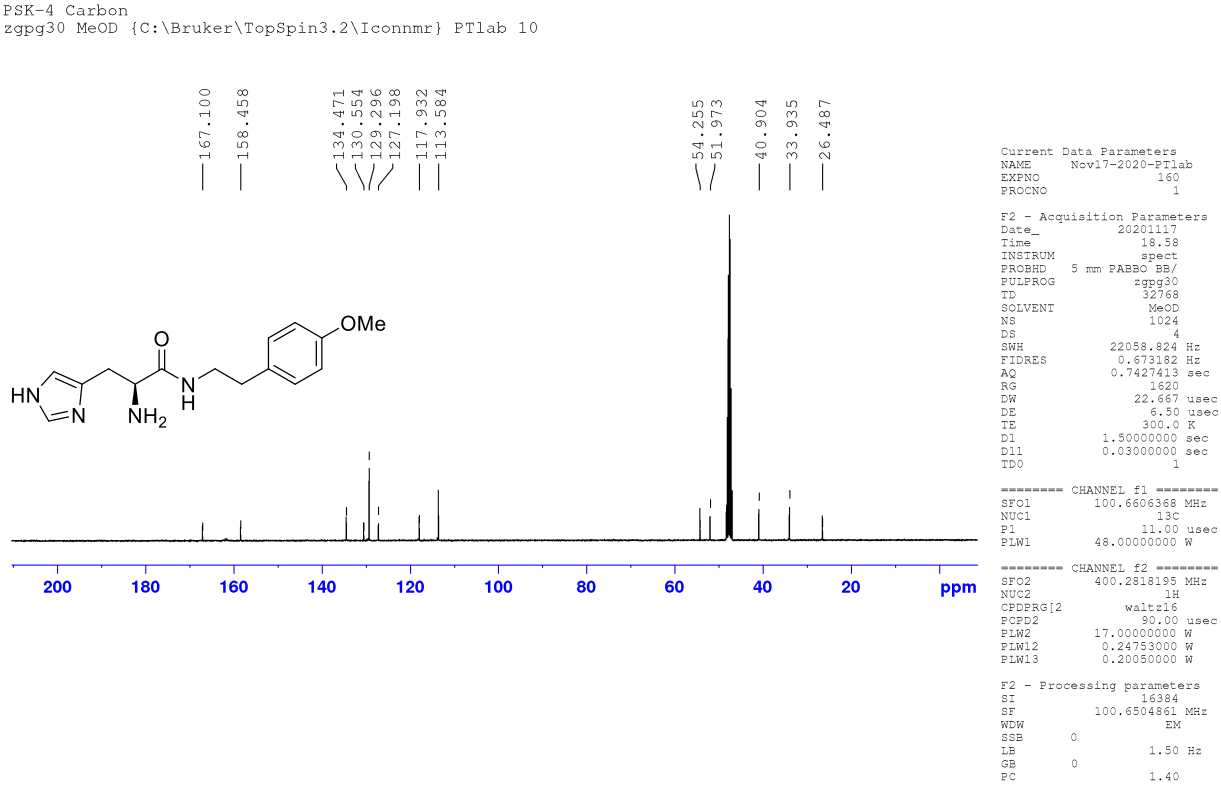
======== CHANNEL f2 ========
SFO2        400.2818195 MHz
NUC2              1H
CPDPRG[2        waltz16
PCPD2           90.00 usec
PLW2        17.00000000 W
PLW12        0.24753000 W
PLW13        0.20050000 W

F2 - Processing parameters
SI                16384
SF          100.6504861 MHz
WDW                EM
SSB       0
LB              1.50 Hz
GB        0
PC              1.40
```

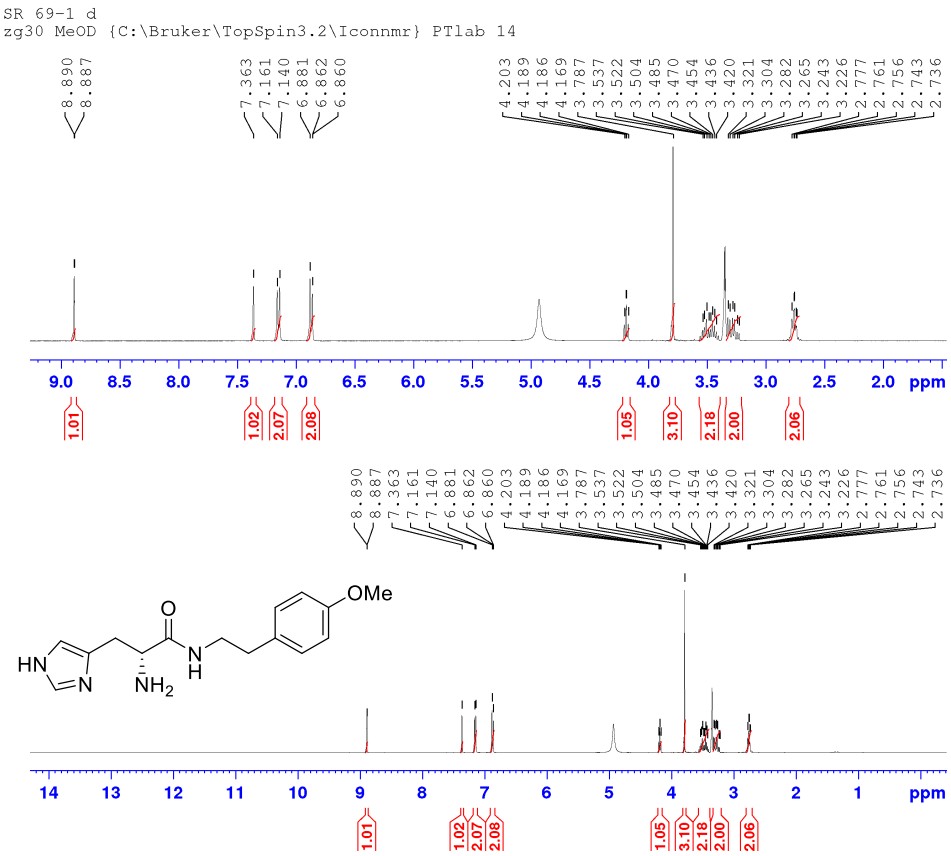
^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-methoxyphenethyl)propanamide **(4e):**



^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-methoxyphenethyl)propanamide **(4e):**



^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-methoxyphenethyl)propanamide **(4f):**

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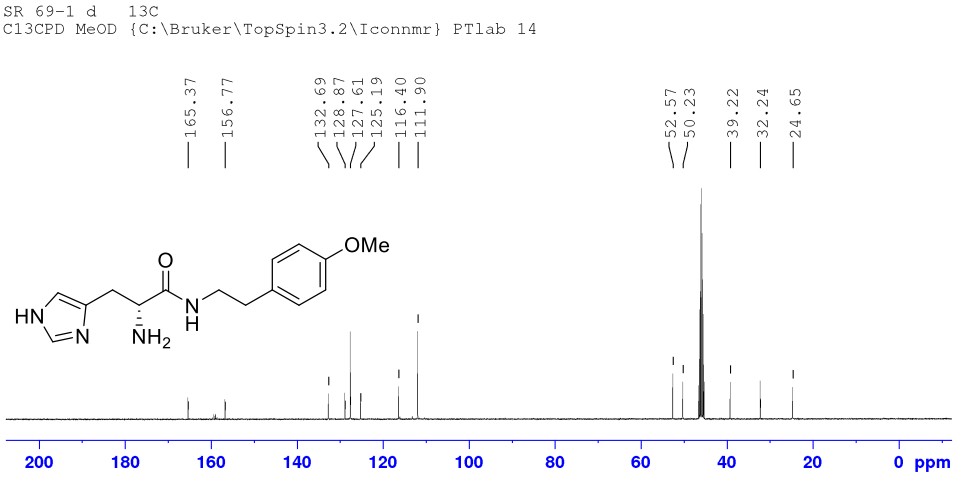
Current Data Parameters
NAME      Nov18-2020-PTlab
EXPNO                  160
PROCNO                   1

F2 - Acquisition Parameters
Date_            20201118
Time               19.34
INSTRUM            spect
PROBHD   5 mm PABBO BB/
PULPROG             zg30
TD                 30046
SOLVENT             MeOD
NS                    16
DS                     1
SWH             8012.820 Hz
FIDRES          0.266685 Hz
AQ              1.8748704 sec
RG                   256
DW                62.400 usec
DE                 6.50 usec
TE                300.0 K
D1             6.00000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1         400.2821952 MHz
NUC1                   1H
P1                 10.86 usec
PLW1         17.00000000 W

F2 - Processing parameters
SI                  131072
SF           400.2793628 MHz
WDW                    EM
SSB          0
LB                  0.20 Hz
GB           0
PC                  1.00
```

^{13}C NMR (100 MHz, MeOH-d_4), (R)-2-amino-3-(1H-imidazol-4-yl)-N-(4-methoxyphenethyl)propanamide **(4f):**

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Current Data Parameters
NAME      Nov18-2020-PTlab
EXPNO                  170
PROCNO                   1

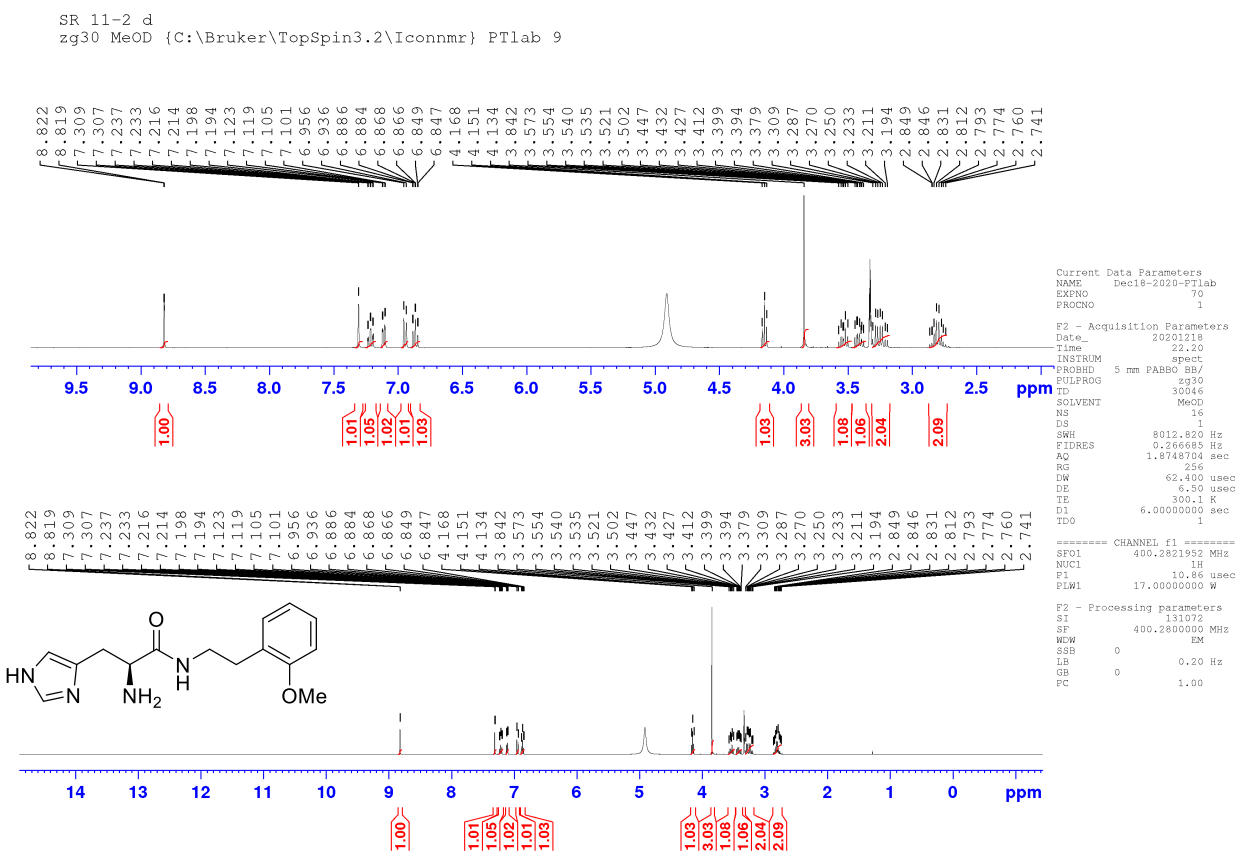
F2 - Acquisition Parameters
Date_           20201118
Time              20.20
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG          zgpg30
TD                65536
SOLVENT            MeOD
NS                  512
DS                    4
SWH           24038.461 Hz
FIDRES         0.366798 Hz
AQ             1.3631488 sec
RG                 2050
DW               20.800 usec
DE                6.50 usec
TE                300.0 K
D1           2.00000000 sec
D11          0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1         100.6605506 MHz
NUC1               13C
P1               11.00 usec
PLW1        48.00000000 W

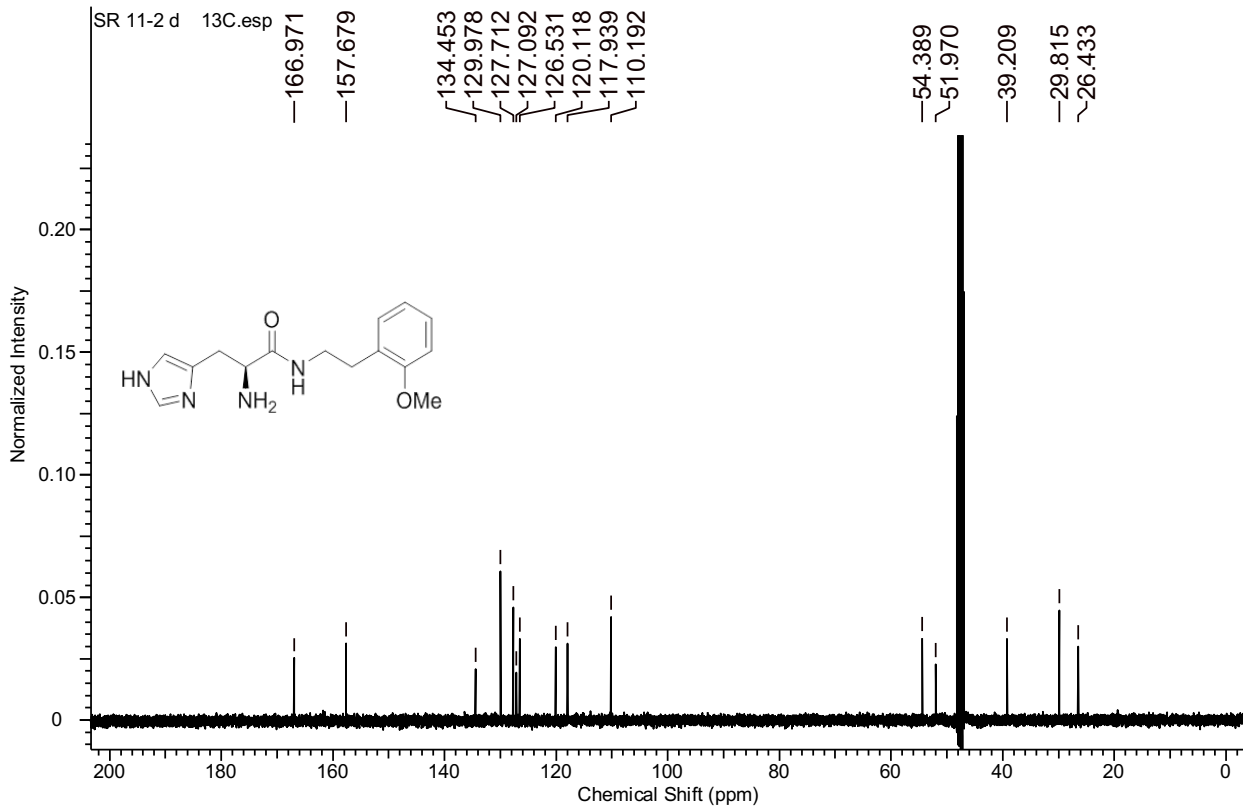
======== CHANNEL f2 ========
SFO2         400.2816011 MHz
NUC2               1H
CPDPRG[2        waltz16
PCPD2            90.00 usec
PLW2        17.00000000 W
PLW12        0.24753000 W
PLW13        0.20050000 W

F2 - Processing parameters
SI                32768
SF           100.6504861 MHz
WDW                EM
SSB       0
LB               1.00 Hz
GB        0
PC               1.40
```

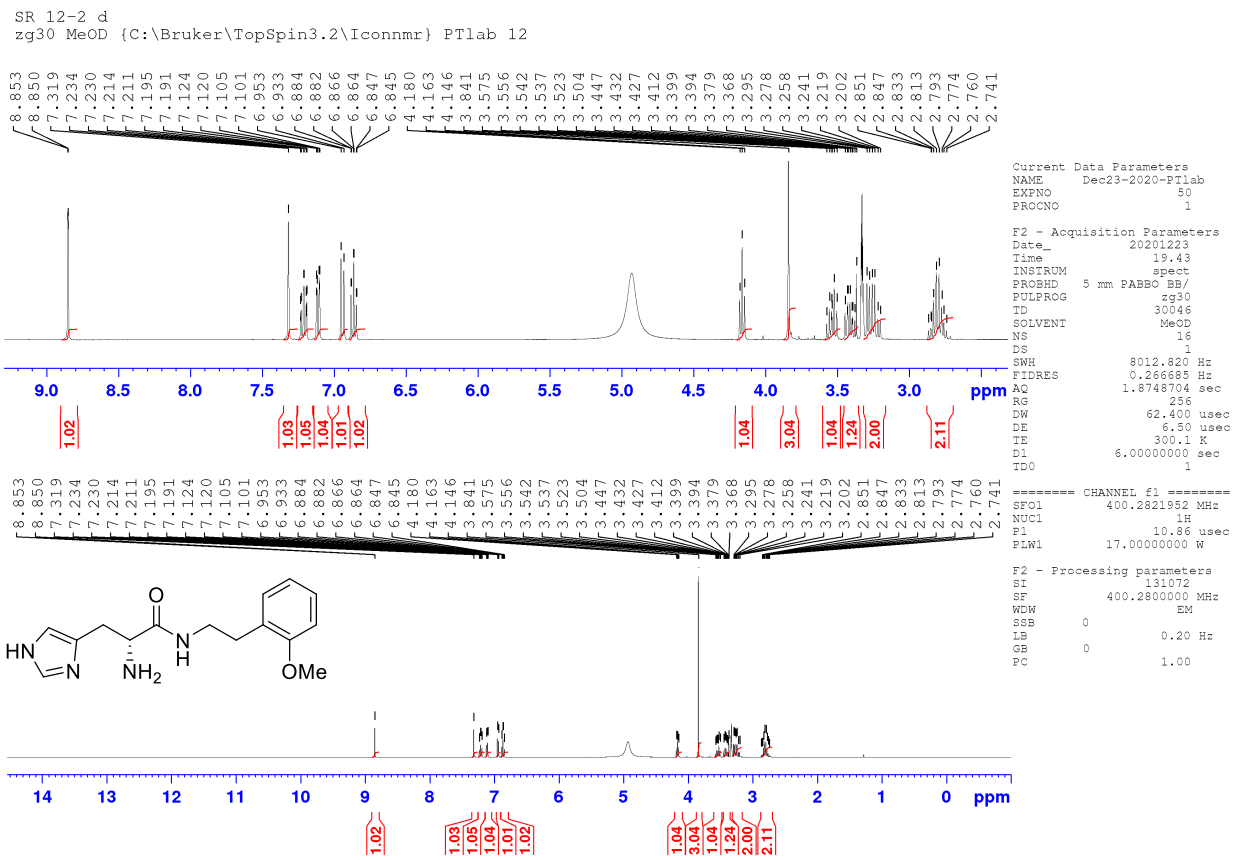
^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-methoxyphenethyl)propanamide **(4g):**



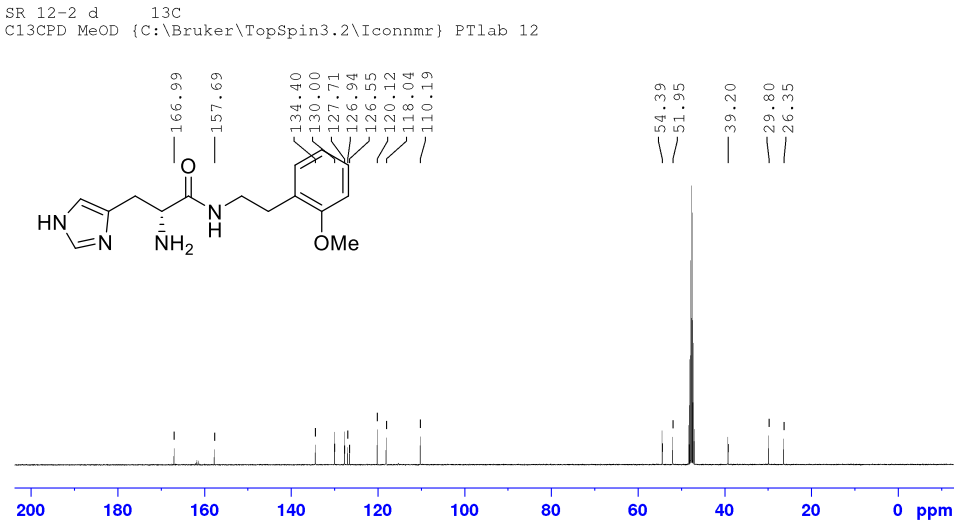
^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-methoxyphenethyl)propanamide **(4g):**



^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-methoxyphenethyl)propanamide (**4h):**



^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-methoxyphenethyl)propanamide (**4h):**

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Current Data Parameters
NAME      Dec23-2020-PTlab
EXPNO                   60
PROCNO                   1

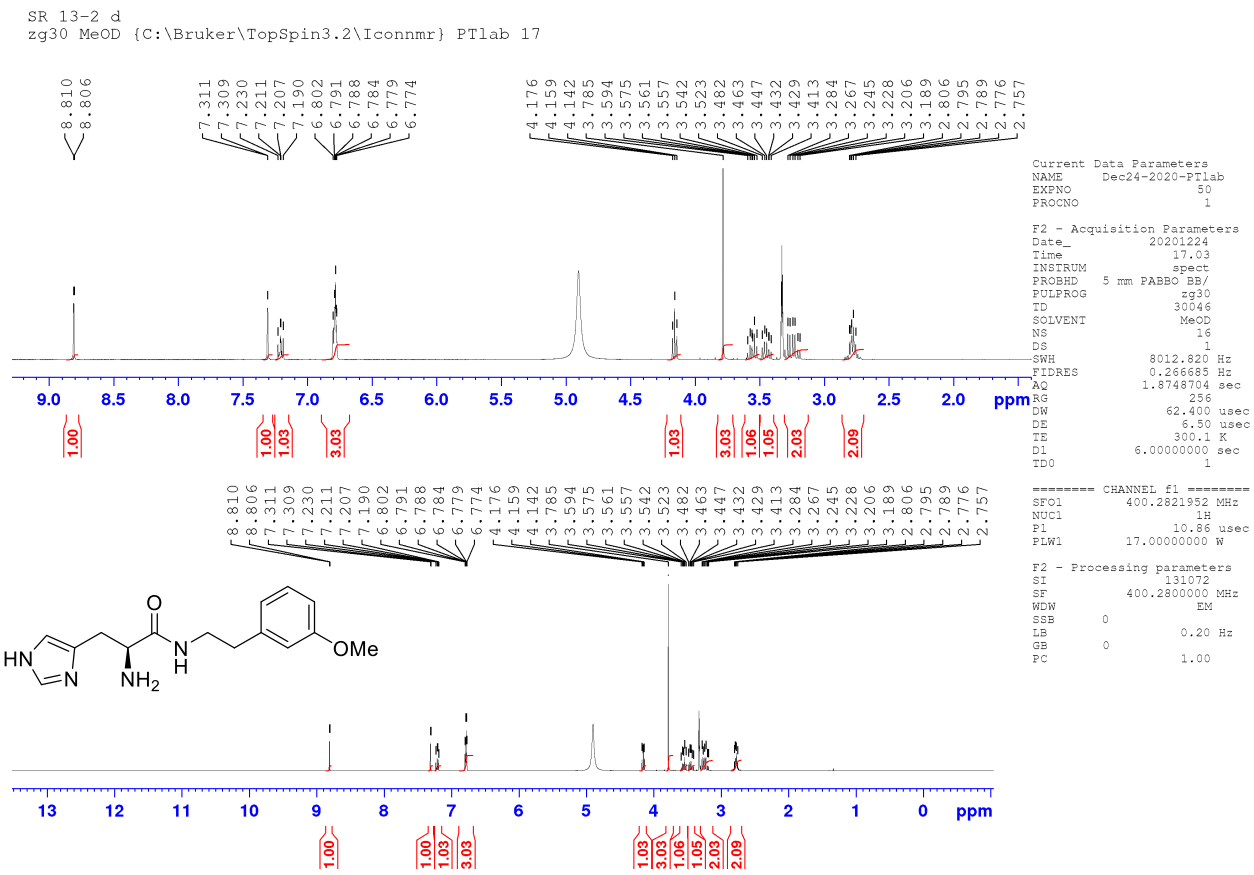
F2 - Acquisition Parameters
Date_              20201223
Time                 20.24
INSTRUM              spect
PROBHD    5 mm PABBO BB/
PULPROG            zgpg30
TD                  65536
SOLVENT               MeOD
NS                     512
DS                       4
SWH              24038.461 Hz
FIDRES            0.366798 Hz
AQ                1.3631488 sec
RG                    2050
DW                  20.800 usec
DE                   6.50 usec
TE                  300.0 K
D1               2.00000000 sec
D11              0.03000000 sec
TD0                      1

======== CHANNEL f1 ========
SFO1          100.6605506 MHz
NUC1                    13C
P1                    11.00 usec
PLW1          48.00000000 W

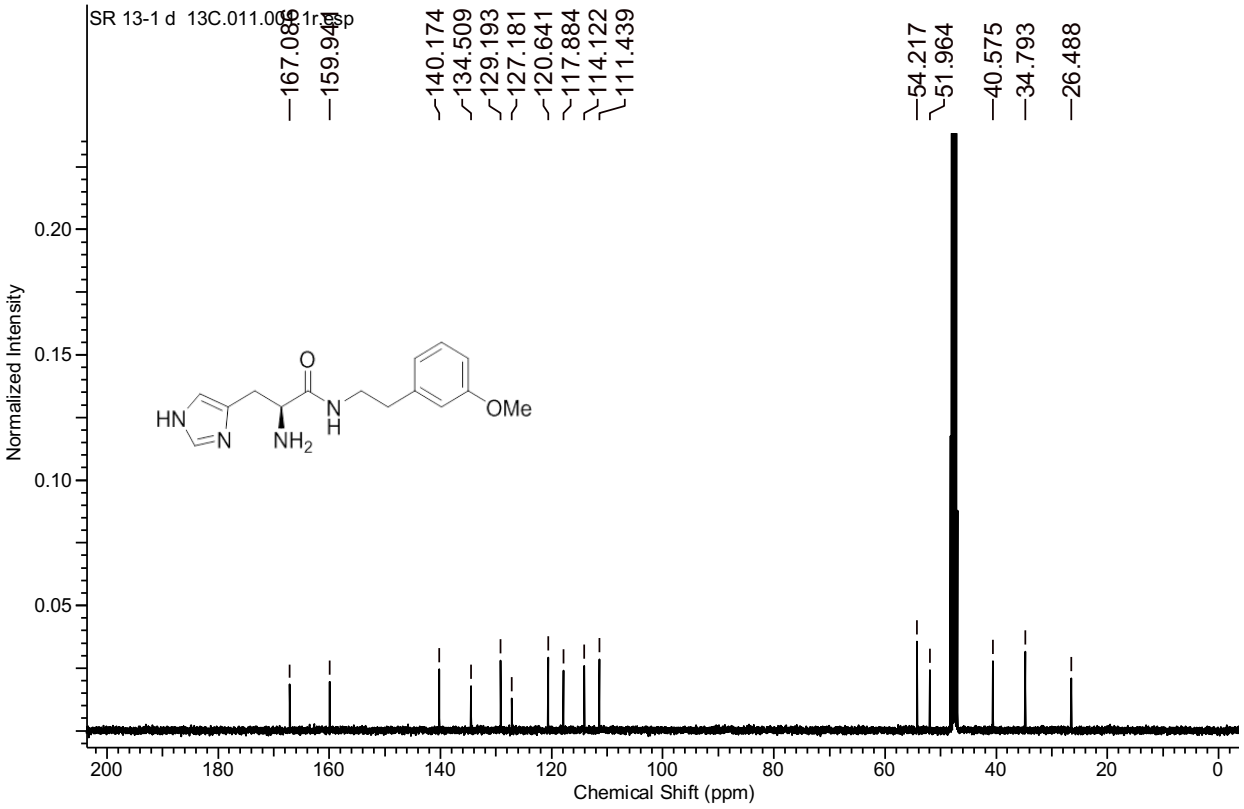
======== CHANNEL f2 ========
SFO2          400.2816011 MHz
NUC2                    1H
CPDPRG[2          waltz16
PCPD2                 90.00 usec
PLW2          17.00000000 W
PLW12          0.24753000 W
PLW13          0.20050000 W

F2 - Processing parameters
SI                    32768
SF            100.6504861 MHz
WDW                      EM
SSB       0
LB                     1.00 Hz
GB        0
PC                     1.40
```

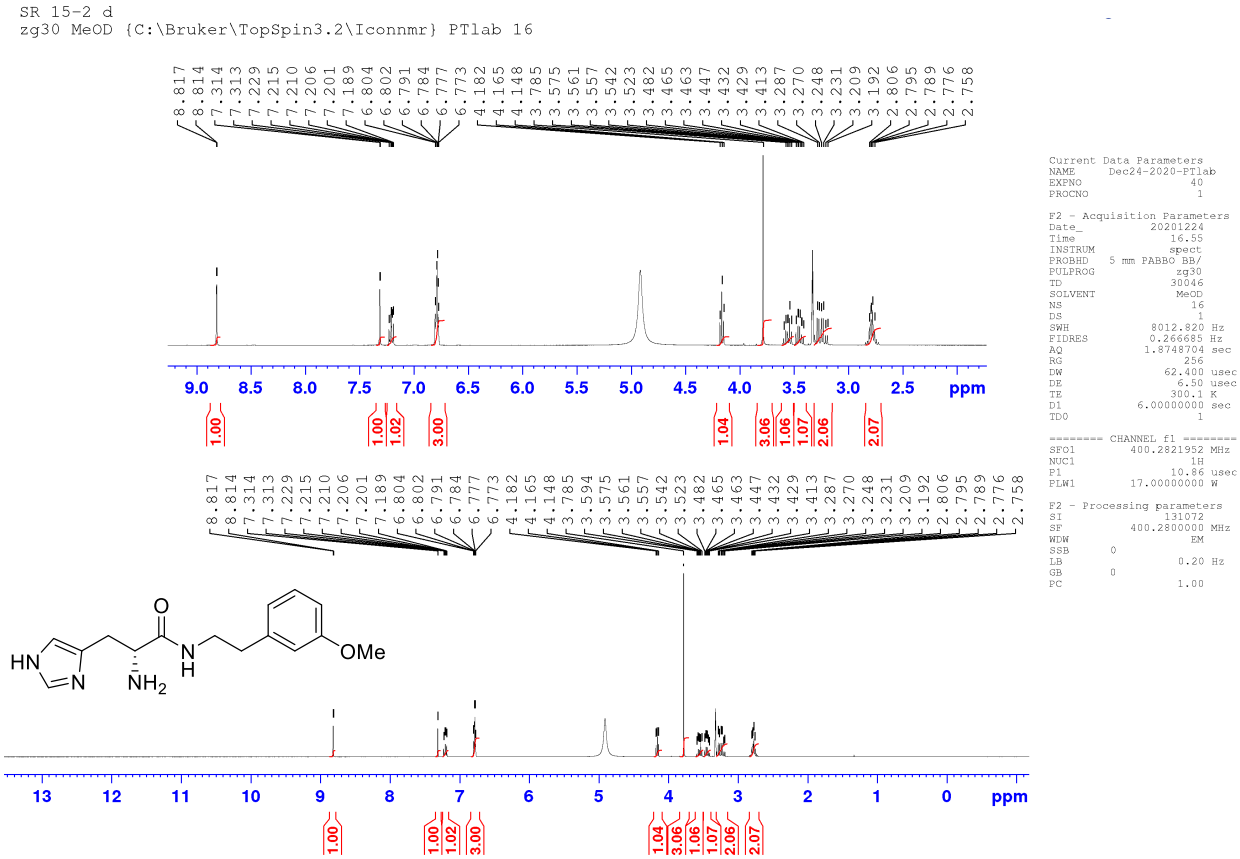
^{1}H NMR (400 MHz, MeOH-d$_4$), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-methoxyphenethyl)propanamide (**4i**)**:**



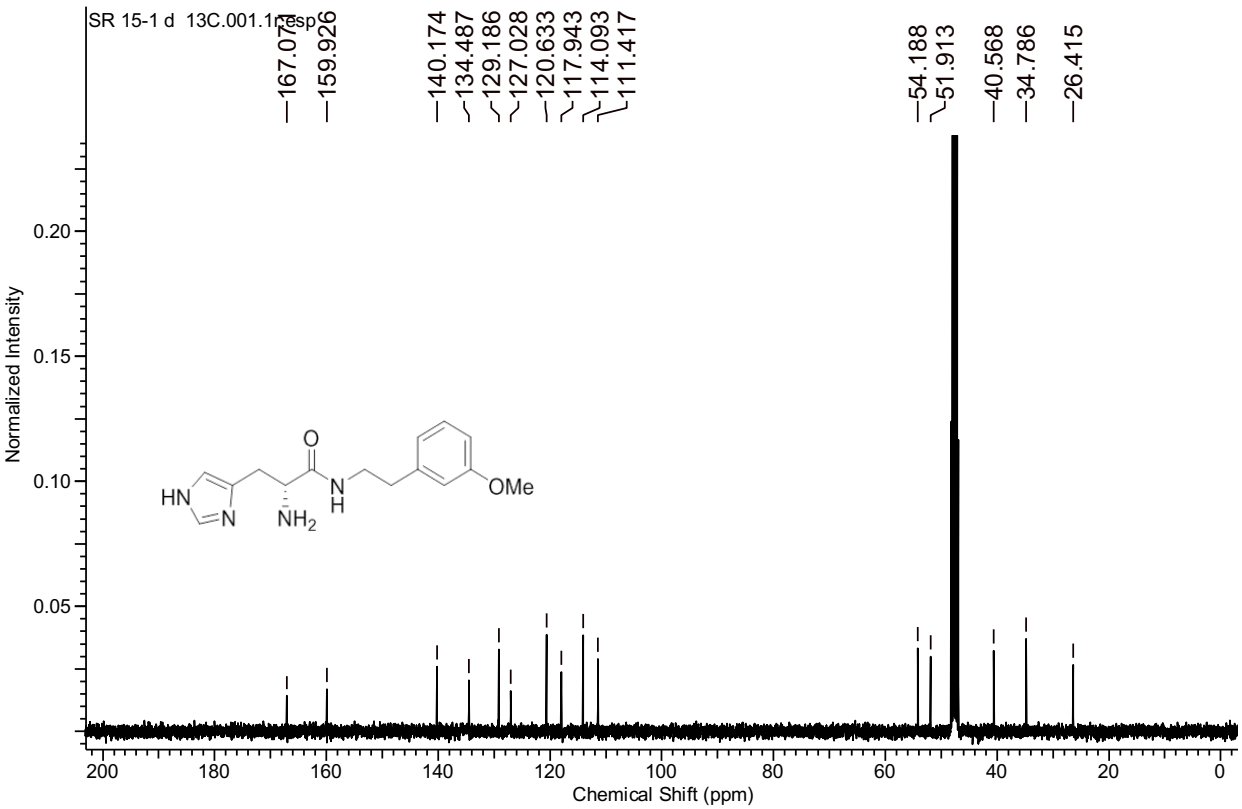
^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-methoxyphenethyl)propanamide (**4i**)**:**



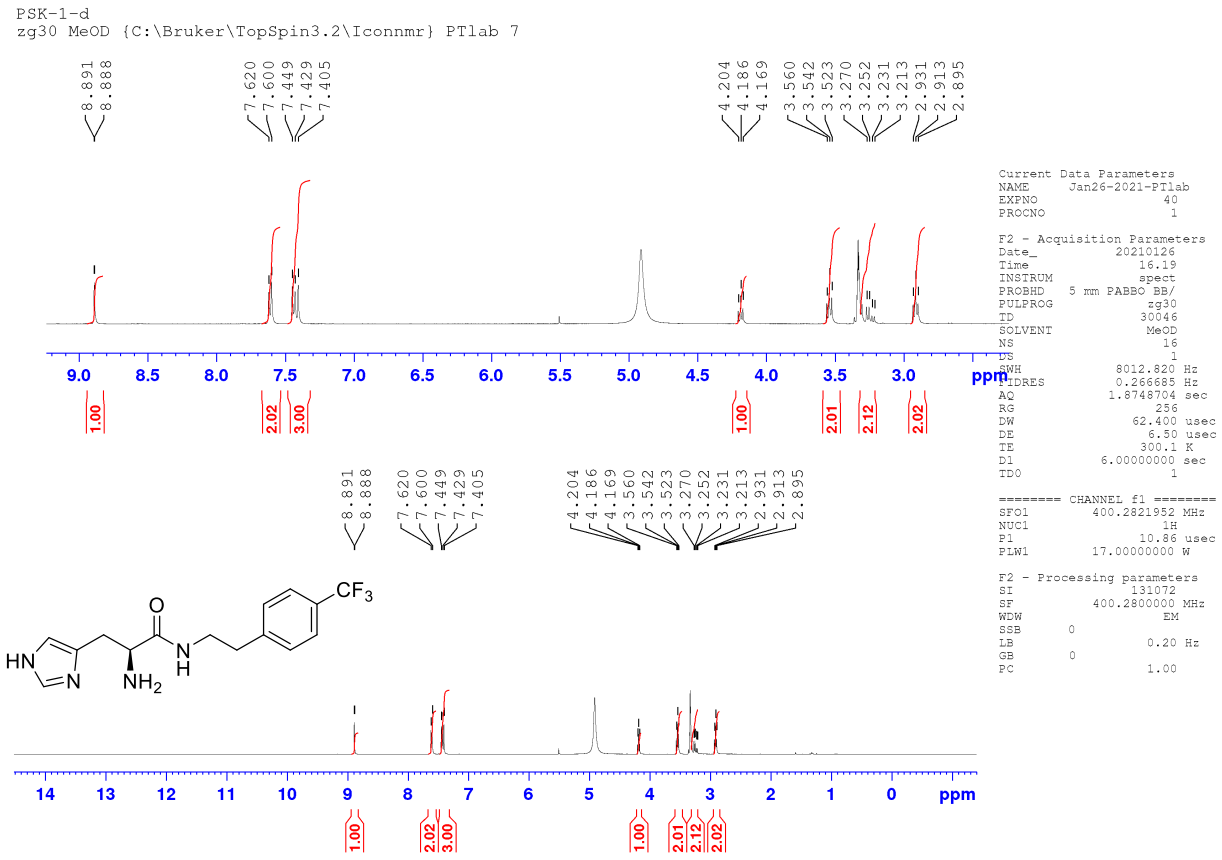
^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-methoxyphenethyl)propanamide (**4j**):



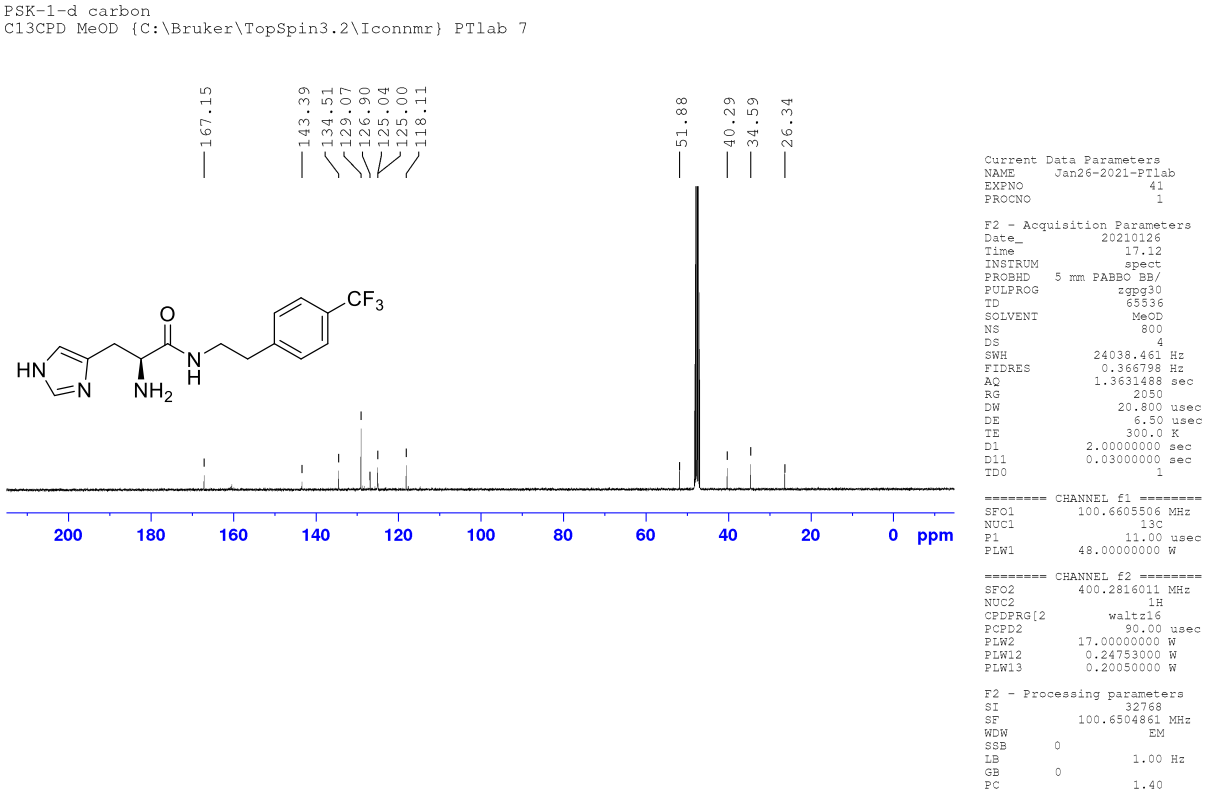
^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-methoxyphenethyl)propanamide (**4j**):



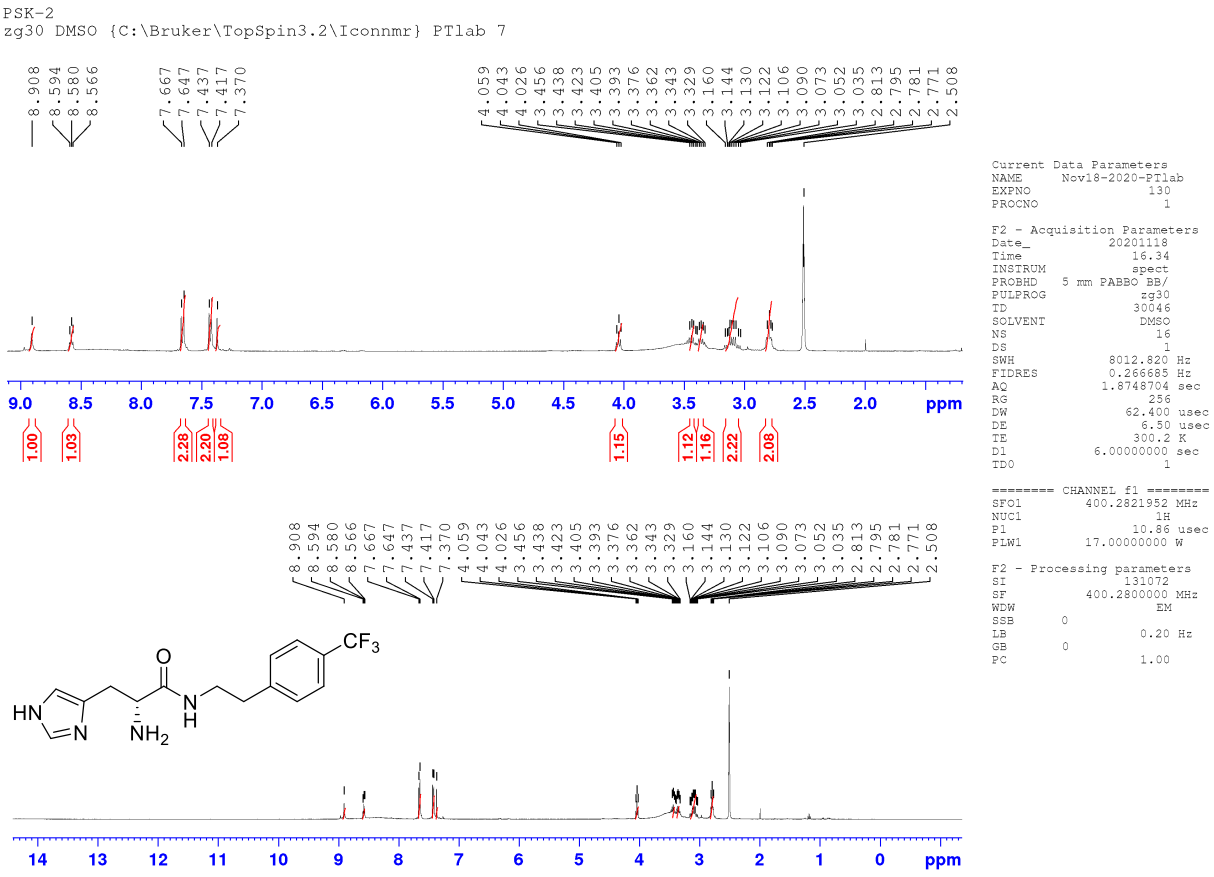
^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4 (trifluoromethyl)phenethyl)propanamide **(4k):**



^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)phenethyl)propanamide **(4k):**



^{1}H NMR (400 MHz, DMSO-d_6), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)phenethyl)propanamide **(4l):**

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EXPNO                  130
PROCNO                   1

F2 - Acquisition Parameters
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Time              16.34
INSTRUM           spect
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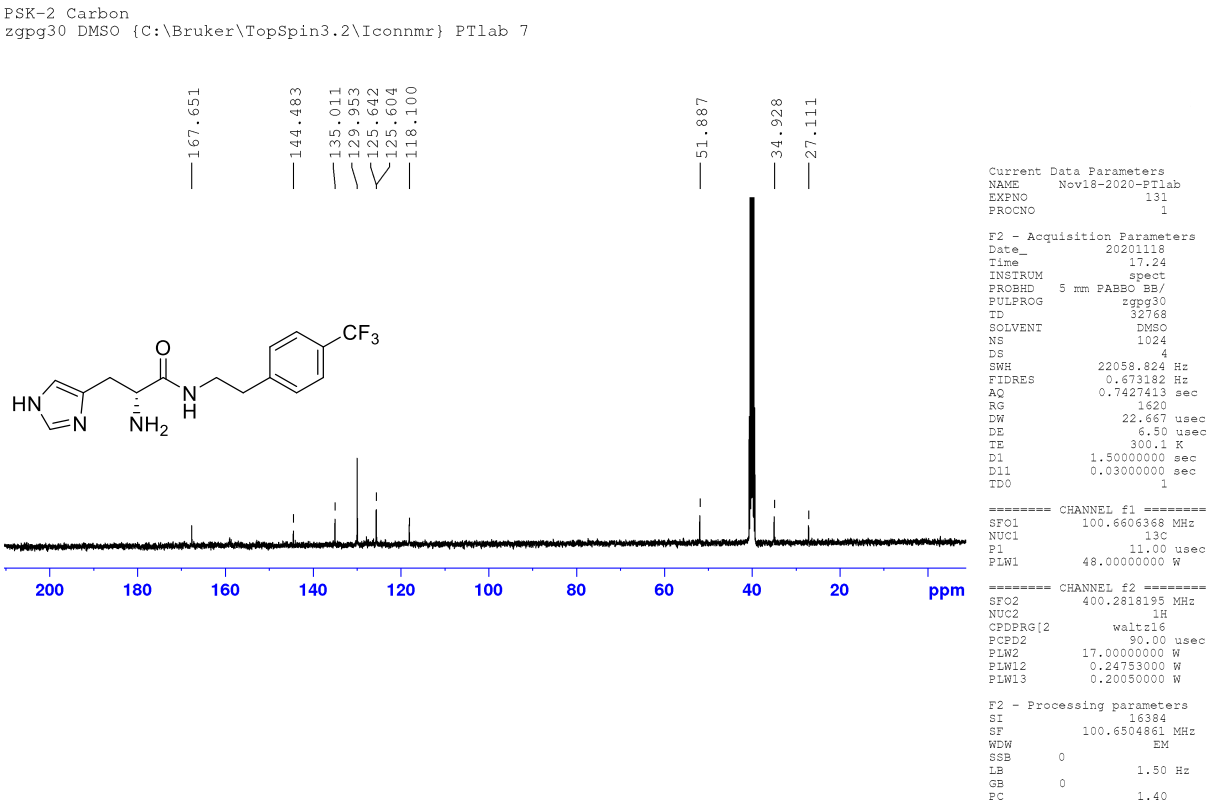
PULPROG             zg30
TD                 30046
SOLVENT             DMSO
NS                    16
DS                     1
SWH            8012.820 Hz
FIDRES         0.266685 Hz
AQ             1.8748704 sec
RG                   256
DW                62.400 usec
DE                 6.50 usec
TE                300.2 K
D1              6.00000000 sec
TD0                    1

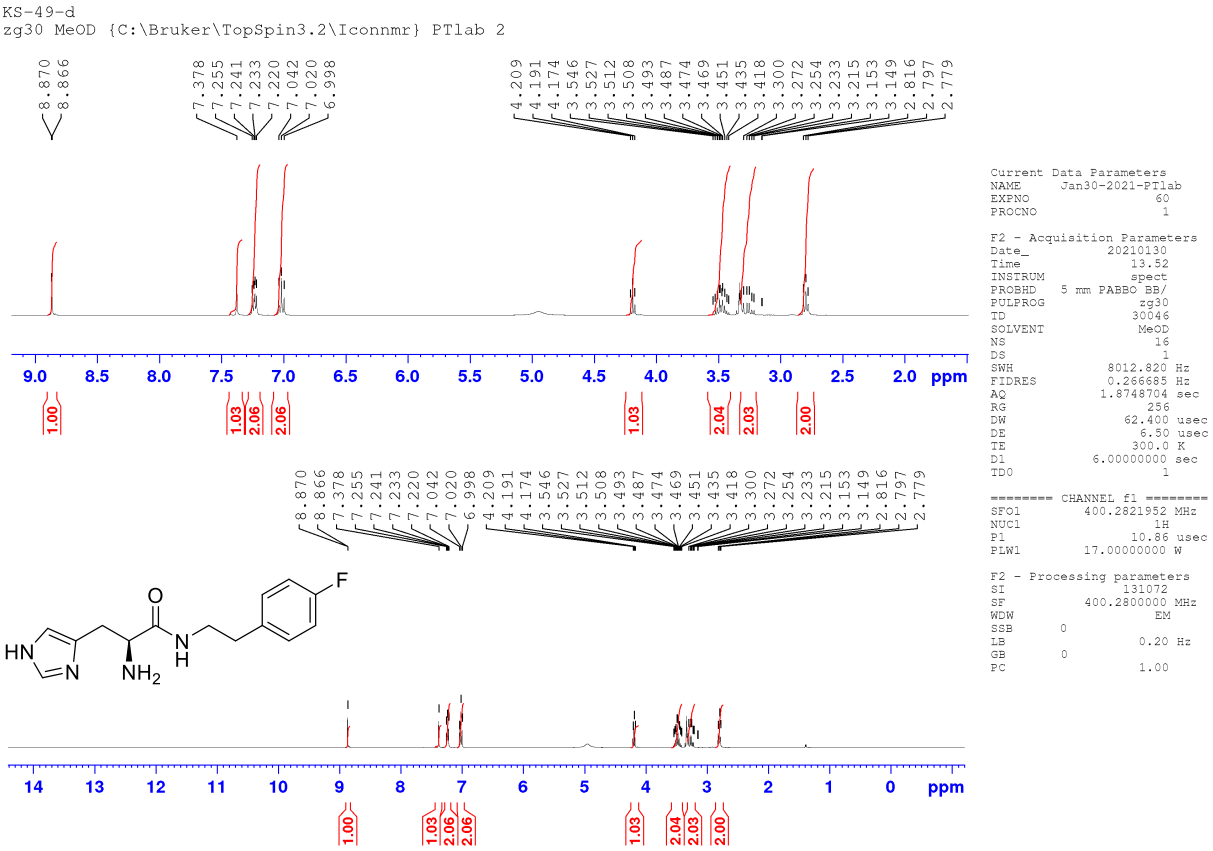
======== CHANNEL f1 ========
SFO1         400.2821952 MHz
NUC1                   1H
P1                 10.86 usec
PLW1        17.00000000 W

F2 - Processing parameters
SI                  131072
SF           400.2800000 MHz
WDW                    EM
SSB       0
LB                0.20 Hz
GB        0
PC                1.00
```

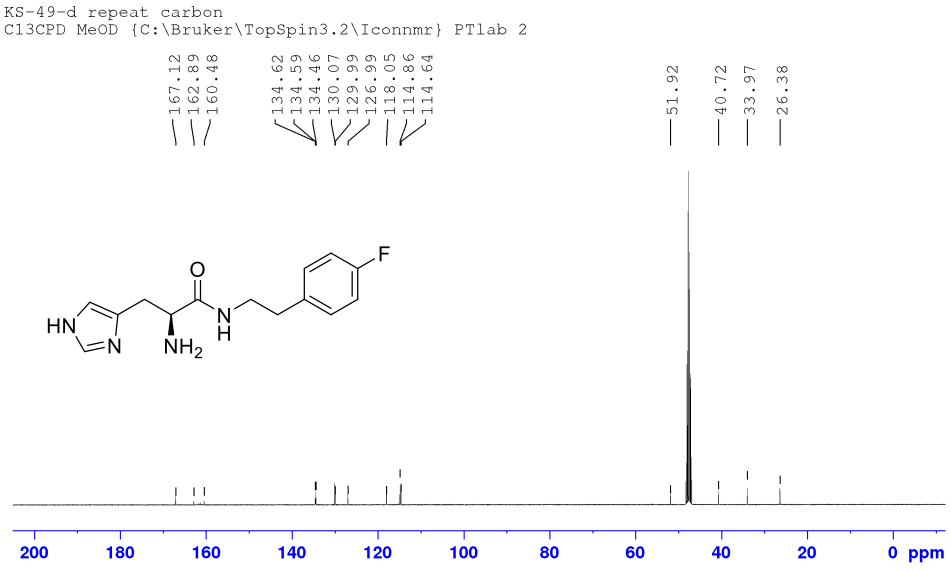
^{13}C NMR (100 MHz, DMSO-d_6), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)phenethyl)propanamide **(4l):**



^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-*N*-(4-fluorophenethyl)-3-(1*H*-imidazol-4-yl)propanamide **(4m):**



^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-*N*-(4-fluorophenethyl)-3-(1*H*-imidazol-4-yl)propanamide **(4m):**

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Current Data Parameters
NAME      Jan30-2021-PTlab
EXPNO                   61
PROCNO                   1

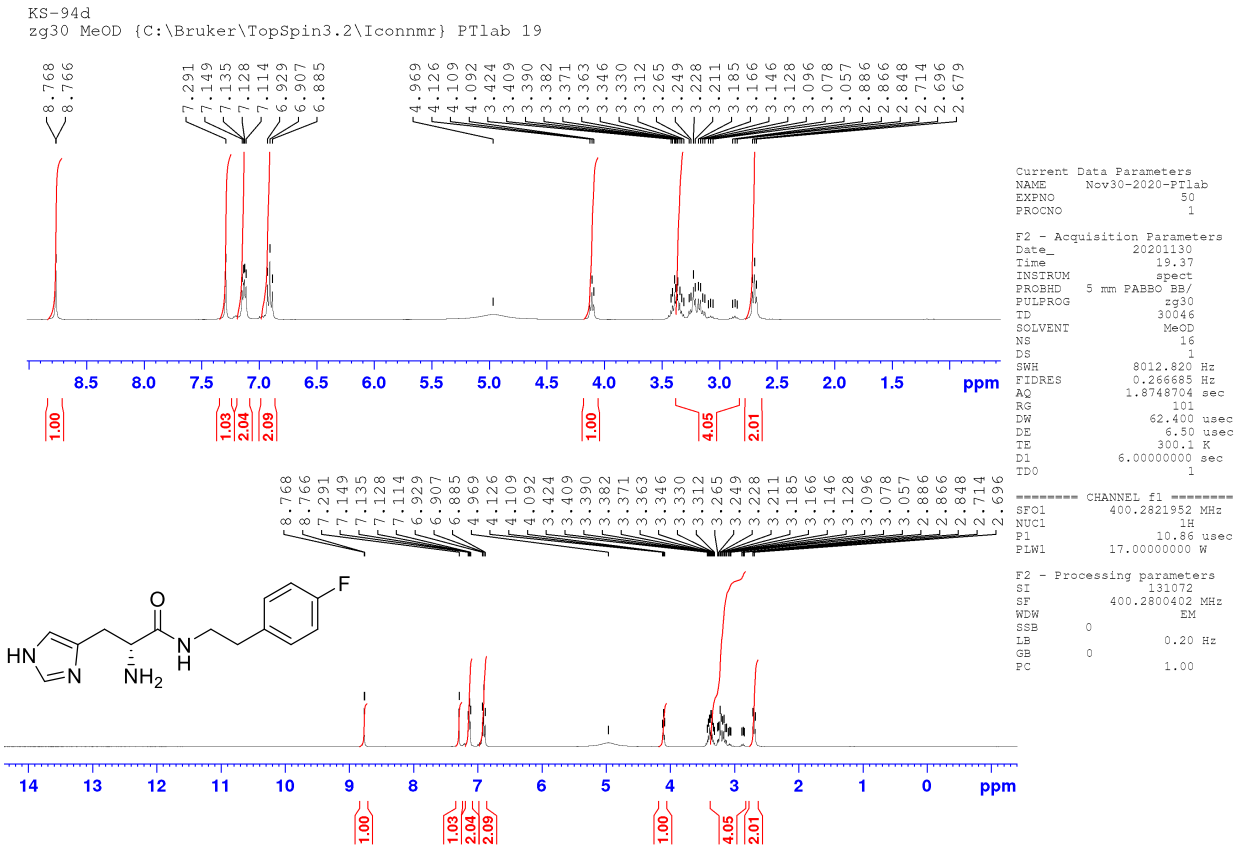
F2 - Acquisition Parameters
Date_           20210130
Time              14.44
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG          zgpg30
TD                65536
SOLVENT             MeOD
NS                    800
DS                      4
SWH           24038.461 Hz
FIDRES         0.366798 Hz
AQ            1.3631488 sec
RG                 2050
DW               20.800 usec
DE                6.50 usec
TE                300.0 K
D1           2.00000000 sec
D11          0.03000000 sec
TD0                     1

======== CHANNEL f1 ========
SFO1         100.6605506 MHz
NUC1                   13C
P1                  11.00 usec
PLW1         48.00000000 W

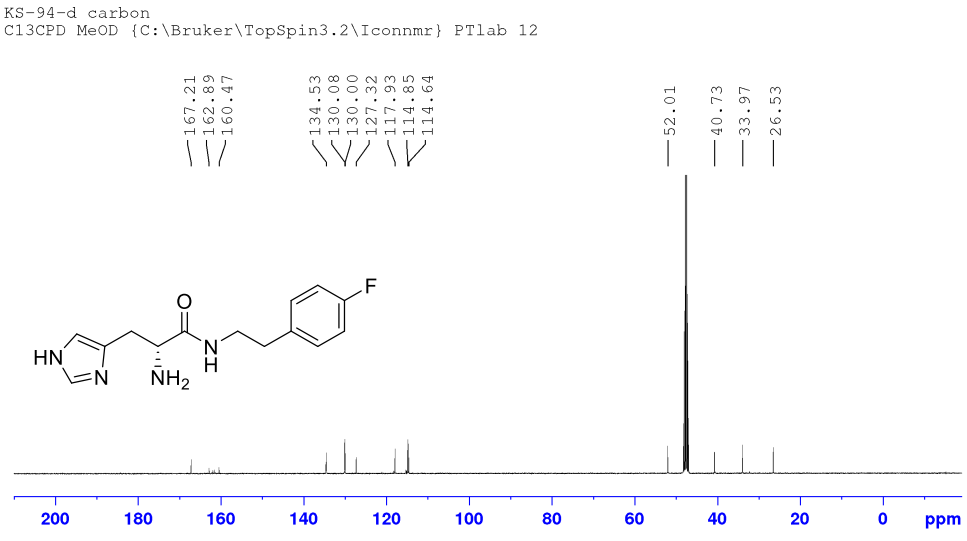
======== CHANNEL f2 ========
SFO2         400.2816011 MHz
NUC2                   1H
CPDPRG[2          waltz16
PCPD2               90.00 usec
PLW2         17.00000000 W
PLW12         0.24753000 W
PLW13         0.20050000 W

F2 - Processing parameters
SI                  32768
SF           100.6504861 MHz
WDW                    EM
SSB       0
LB                  1.00 Hz
GB        0
PC                  1.40
```

^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-*N*-(4-fluorophenethyl)-3-(1*H*-imidazol-4-yl)propanamide **(4n):**



^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-*N*-(4-fluorophenethyl)-3-(1*H*-imidazol-4-yl)propanamide **(4n):**

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Current Data Parameters
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EXPNO                  111
PROCNO                   1

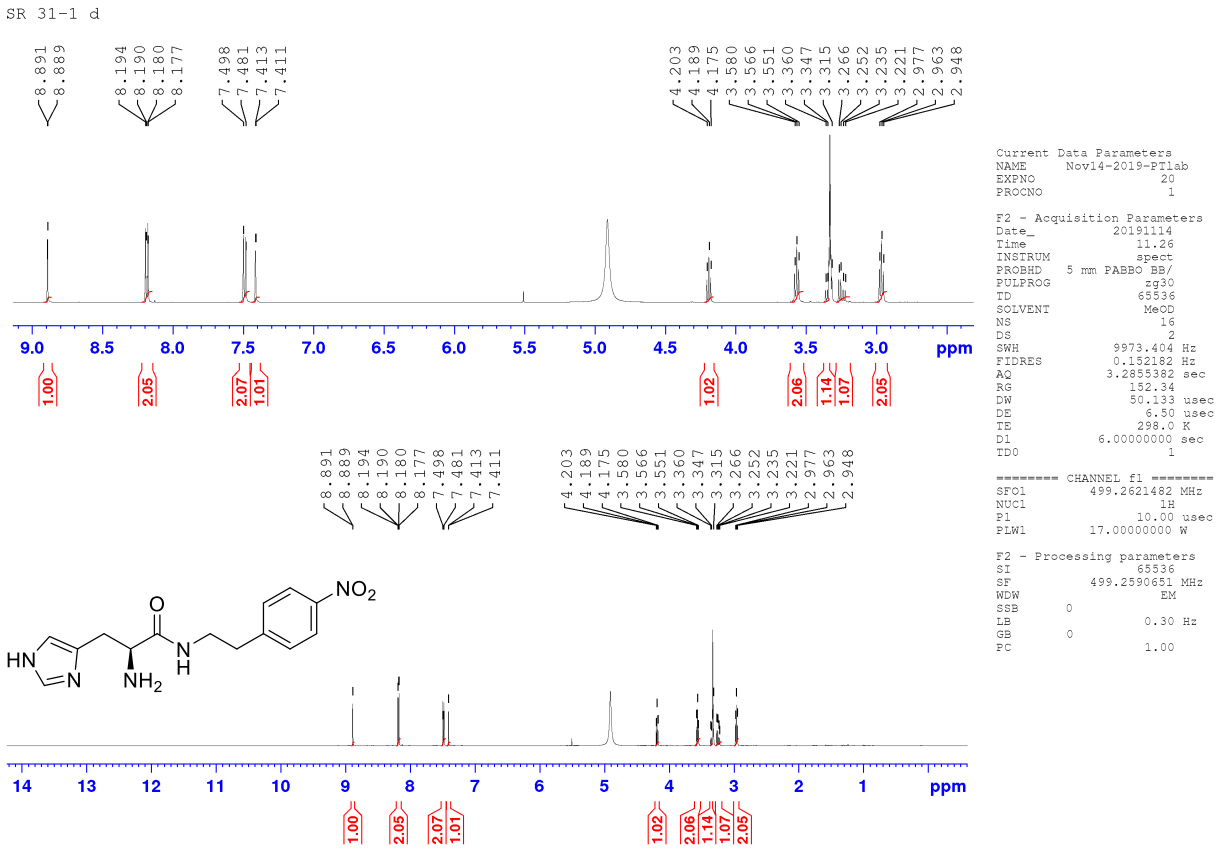
F2 - Acquisition Parameters
Date_           20210129
Time              18.10
INSTRUM             spect
PROBHD    5 mm PABBO BB/
PULPROG           zgpg30
TD                65536
SOLVENT             MeOD
NS                  800
DS                    4
SWH           24038.461 Hz
FIDRES         0.366798 Hz
AQ             1.3631488 sec
RG                 2050
DW                20.800 usec
DE                 6.50 usec
TE                300.1 K
D1            2.00000000 sec
D11           0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1         100.6605506 MHz
NUC1                 13C
P1                 11.00 usec
PLW1        48.00000000 W

======== CHANNEL f2 ========
SFO2         400.2816011 MHz
NUC2                 1H
CPDPRG[2          waltz16
PCPD2              90.00 usec
PLW2        17.00000000 W
PLW12        0.24753000 W
PLW13        0.20050000 W

F2 - Processing parameters
SI                 32768
SF           100.6504861 MHz
WDW                  EM
SSB       0
LB                 1.00 Hz
GB        0
PC                 1.40
```

^{1}H NMR (500 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-nitrophenethyl)propanamide (**4o**):



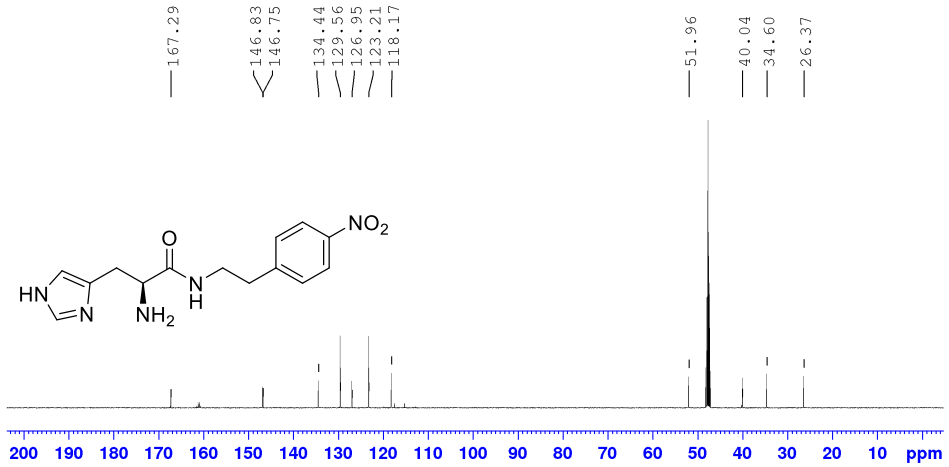

```
Current Data Parameters
NAME      Nov14-2019-PTlab
EXPNO                   20
PROCNO                   1

F2 - Acquisition Parameters
Date_            20191114
Time                11.26
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG             zg30
TD                 65536
SOLVENT             MeOD
NS                    16
DS                     2
SWH              9973.404 Hz
FIDRES           0.152182 Hz
AQ               3.2855382 sec
RG                 152.34
DW                50.133 usec
DE                 6.50 usec
TE                298.0 K
D1             6.00000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1         499.2621482 MHz
NUC1                   1H
P1                 10.00 usec
PLW1         17.00000000 W

F2 - Processing parameters
SI                 65536
SF           499.2590651 MHz
WDW                   EM
SSB       0
LB                 0.30 Hz
GB        0
PC                 1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-nitrophenethyl)propanamide (**4o**):




```
Current Data Parameters
NAME      Nov15-2019-PTlab
EXPNO                   40
PROCNO                   1

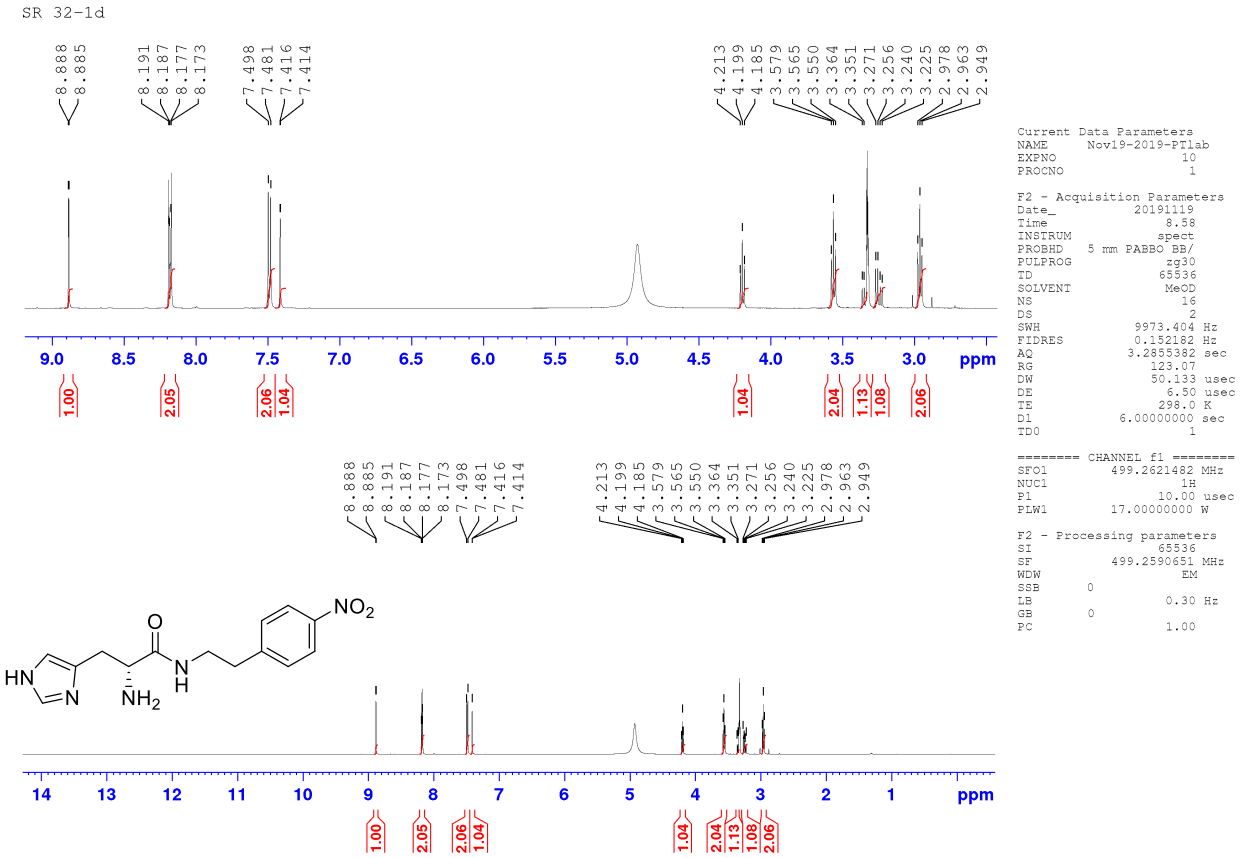
F2 - Acquisition Parameters
Date_           20191115
Time              20.25
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG          zgpg30
TD                65536
SOLVENT            MeOD
NS                  512
DS                    4
SWH           29761.904 Hz
FIDRES         0.454131 Hz
AQ            1.1010048 sec
RG               191.52
DW               16.800 usec
DE               10.00 usec
TE               300.0 K
D1           2.00000000 sec
D11          0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1        125.5513451 MHz
NUC1               13C
P1               10.00 usec
PLW1        71.00000000 W

======== CHANNEL f2 ========
SFO2        499.2610621 MHz
NUC2               1H
CPDPRG[2         waltz16
PCPD2             80.00 usec
PLW2        17.00000000 W
PLW12        0.26563001 W
PLW13        0.17000000 W

F2 - Processing parameters
SI                32768
SF          125.5387918 MHz
WDW                EM
SSB       0
LB               1.00 Hz
GB        0
PC               1.40
```

^{1}H NMR (500 MHz, MeOH-d_4), (R)-2-amino-3-(1H-imidazol-4-yl)-N-(4-nitrophenethyl)propanamide (**4p**):



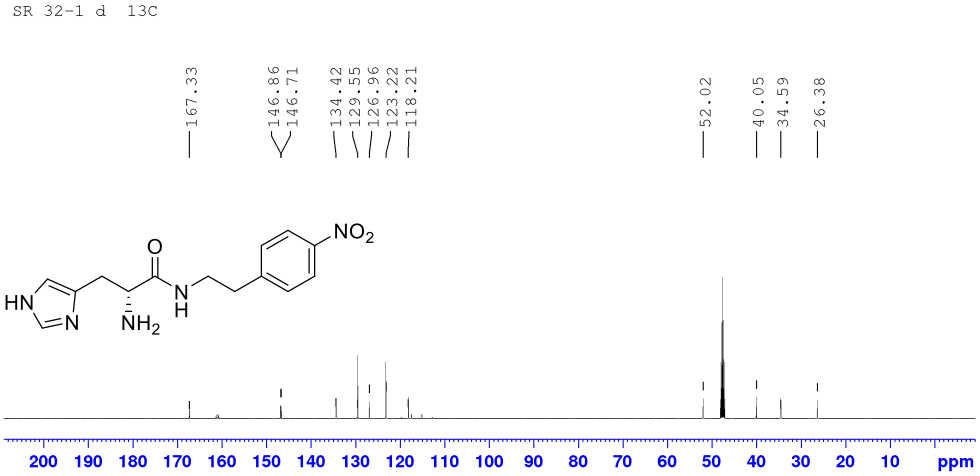

```
Current Data Parameters
NAME      Nov19-2019-PTlab
EXPNO                   10
PROCNO                   1

F2 - Acquisition Parameters
Date_            20191119
Time                 8.58
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG             zg30
TD                 65536
SOLVENT             MeOD
NS                    16
DS                     2
SWH             9973.404 Hz
FIDRES          0.152182 Hz
AQ              3.2855382 sec
RG                 123.07
DW                50.133 usec
DE                 6.50 usec
TE                298.0 K
D1             6.00000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1          499.2621482 MHz
NUC1                   1H
P1                 10.00 usec
PLW1          17.00000000 W

F2 - Processing parameters
SI                 65536
SF            499.2590651 MHz
WDW                   EM
SSB       0
LB                 0.30 Hz
GB        0
PC                 1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-nitrophenethyl)propanamide (**4p):**




```
Current Data Parameters
NAME       Nov20-2019-PTlab
EXPNO                   50
PROCNO                   1

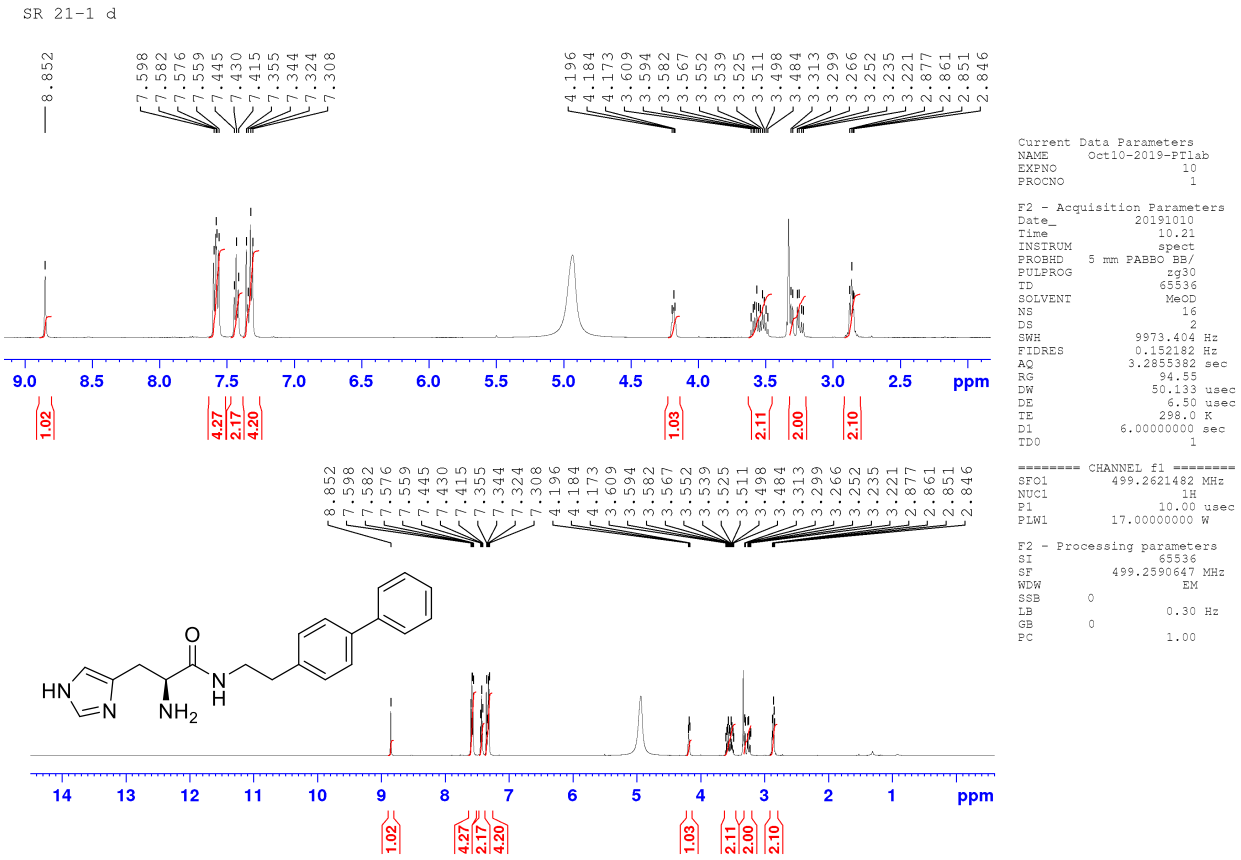
F2 - Acquisition Parameters
Date_            20191120
Time                17.06
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG          zgpg30
TD                65536
SOLVENT            MeOD
NS                  512
DS                    4
SWH           29761.904 Hz
FIDRES         0.454131 Hz
AQ             1.1010048 sec
RG                191.52
DW               16.800 usec
DE               10.00 usec
TE                300.0 K
D1           2.00000000 sec
D11          0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1        125.5513451 MHz
NUC1                  13C
P1                  10.00 usec
PLW1        71.00000000 W

======== CHANNEL f2 ========
SFO2        499.2610621 MHz
NUC2                  1H
CPDPRG[2         waltz16
PCPD2               80.00 usec
PLW2        17.00000000 W
PLW12        0.26563001 W
PLW13        0.17000000 W

F2 - Processing parameters
SI                  32768
SF          125.5387918 MHz
WDW                   EM
SSB       0
LB                  1.00 Hz
GB        0
PC                  1.40
```

^{1}H NMR (500 MHz, MeOH-d_4), (*S*)-*N*-(2-([1,1'-biphenyl]-4-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide (**4q**):



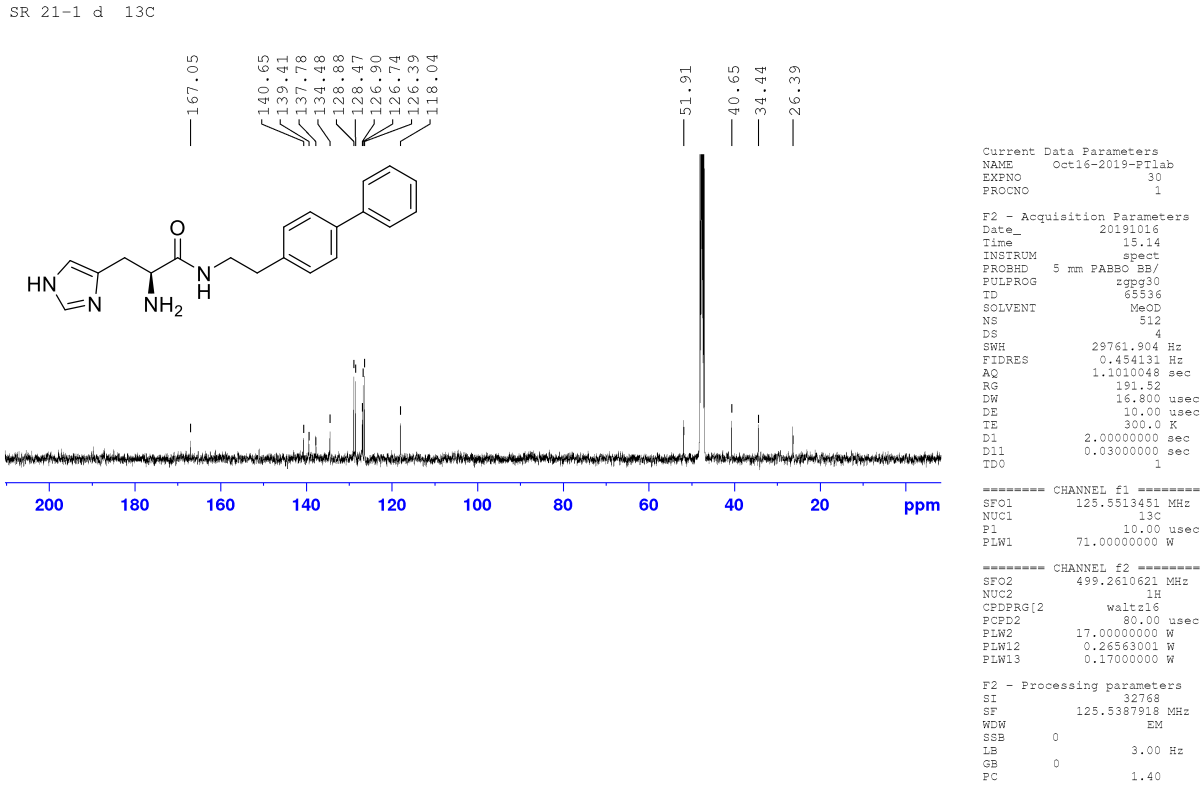

```
Current Data Parameters
NAME      Oct10-2019-Ptlab
EXPNO                  10
PROCNO                  1

F2 - Acquisition Parameters
Date_              20191010
Time                 10.21
INSTRUM              spect
PROBHD   5 mm PABBO BB/
PULPROG              zg30
TD                  65536
SOLVENT              MeOD
NS                     16
DS                      2
SWH              9973.404 Hz
FIDRES           0.152182 Hz
AQ               3.2855382 sec
RG                  94.55
DW                 50.133 usec
DE                  6.50 usec
TE                 298.0 K
D1              6.00000000 sec
TD0                     1

======== CHANNEL f1 ========
SFO1         499.2621482 MHz
NUC1                   1H
P1                  10.00 usec
PLW1         17.00000000 W

F2 - Processing parameters
SI                  65536
SF           499.2590647 MHz
WDW                    EM
SSB      0
LB                  0.30 Hz
GB       0
PC                  1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*S*)-*N*-(2-([1,1'-biphenyl]-4-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide (**4q**):




```
Current Data Parameters
NAME      Oct16-2019-PTlab
EXPNO                   30
PROCNO                   1

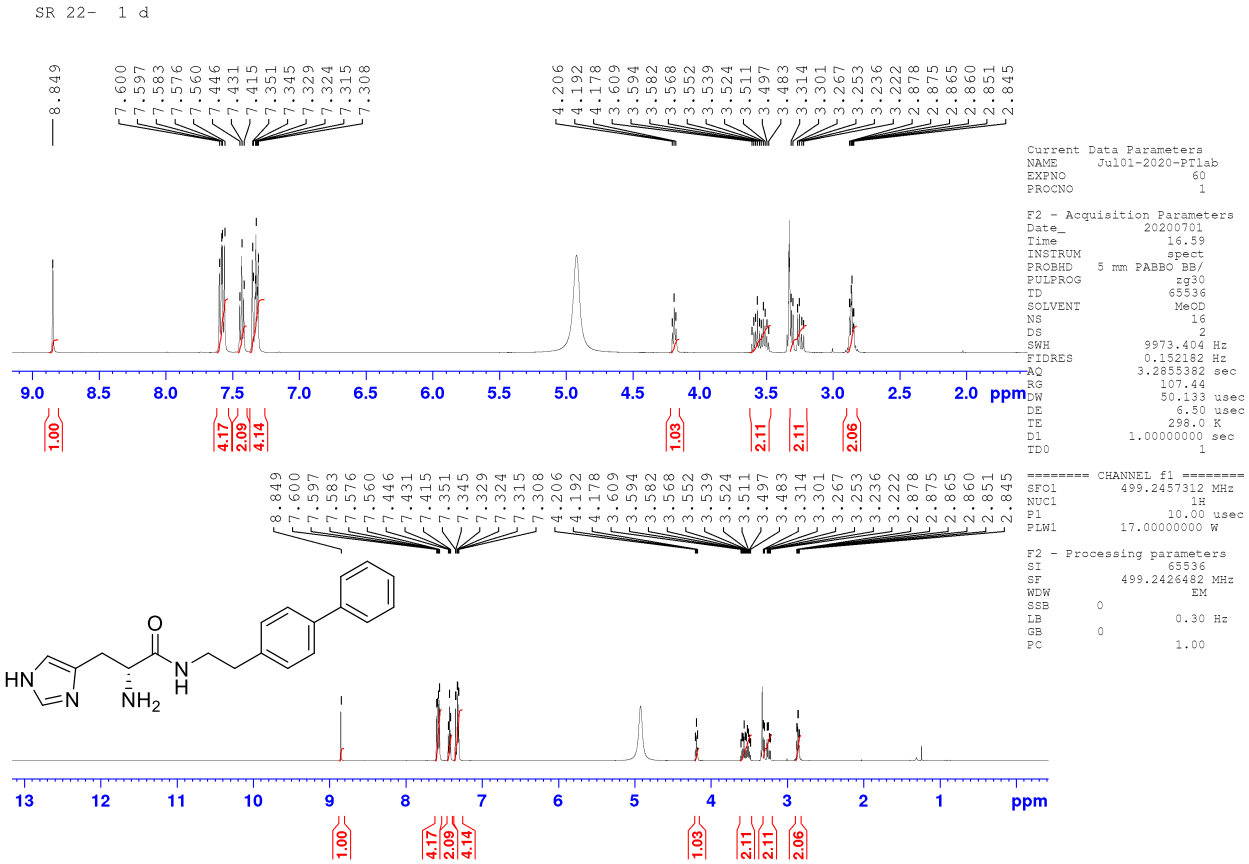
F2 - Acquisition Parameters
Date_           20191016
Time              15.14
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG          zgpg30
TD                65536
SOLVENT             MeOD
NS                    512
DS                      4
SWH           29761.904 Hz
FIDRES         0.454131 Hz
AQ            1.1010048 sec
RG               191.52
DW               16.800 usec
DE               10.00 usec
TE               300.0 K
D1            2.00000000 sec
D11           0.03000000 sec
TD0                     1

======== CHANNEL f1 ========
SFO1         125.5513451 MHz
NUC1               13C
P1               10.00 usec
PLW1         71.00000000 W

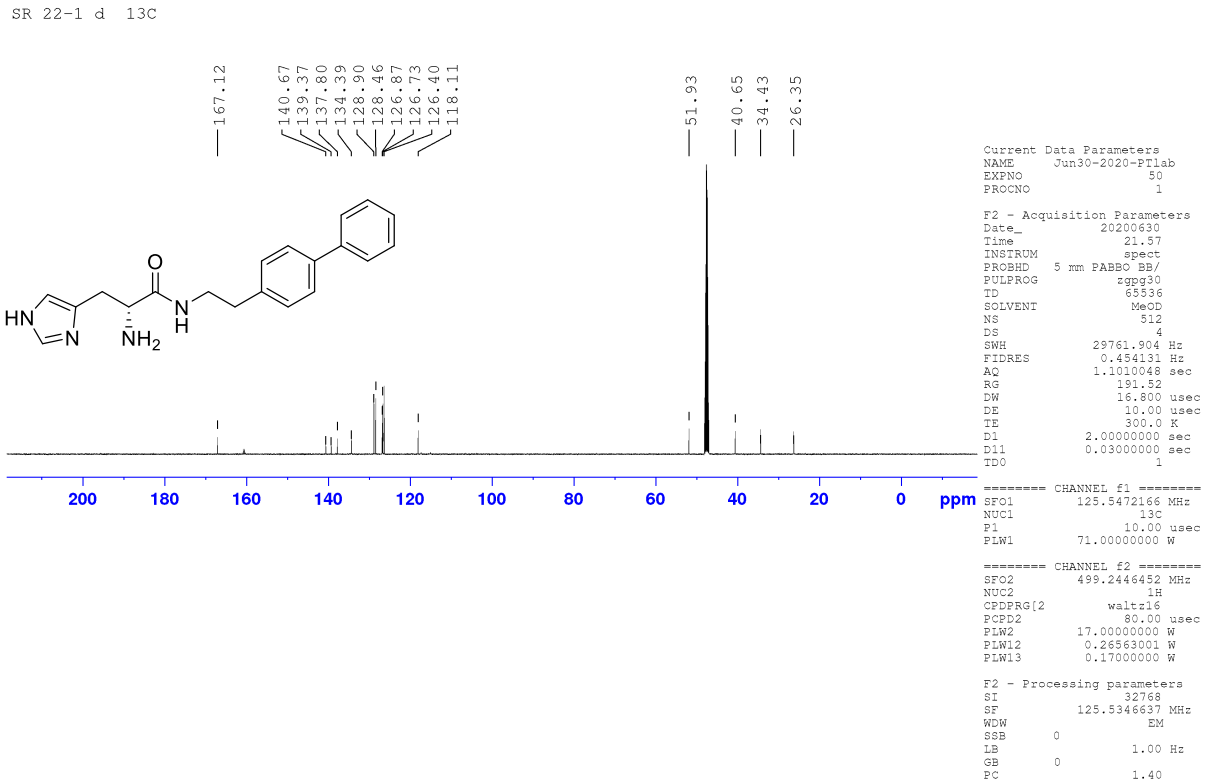
======== CHANNEL f2 ========
SFO2         499.2610621 MHz
NUC2               1H
CPDPRG[2         waltz16
PCPD2             80.00 usec
PLW2         17.00000000 W
PLW12         0.26563001 W
PLW13         0.17000000 W

F2 - Processing parameters
SI                 32768
SF           125.5387918 MHz
WDW                  EM
SSB       0
LB                3.00 Hz
GB        0
PC                1.40
```

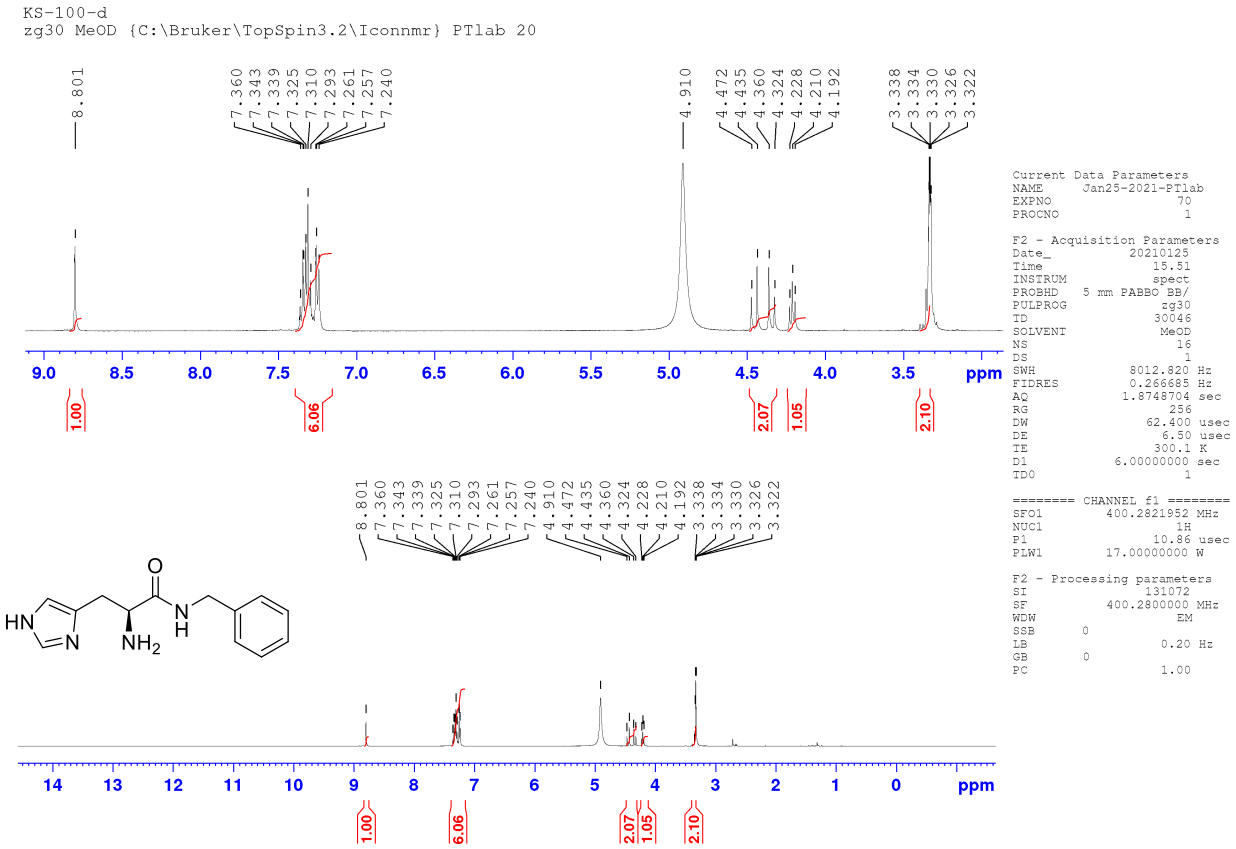
^{1}H NMR (500 MHz, MeOH-d_4), (*R*)-*N*-(2-([1,1'-biphenyl]-4-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide (**4r**):



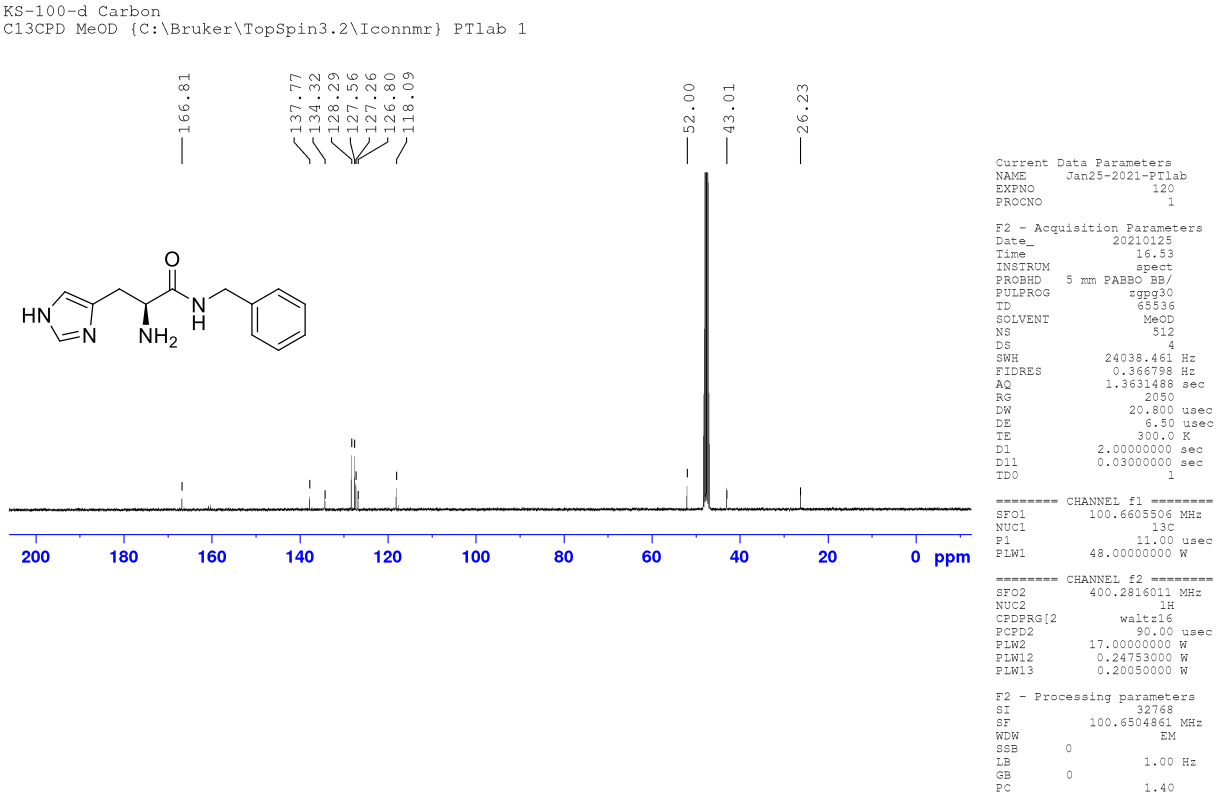
^{13}C NMR (126 MHz, MeOH-d_4), (*R*)-*N*-(2-([1,1'-biphenyl]-4-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide (**4r**):



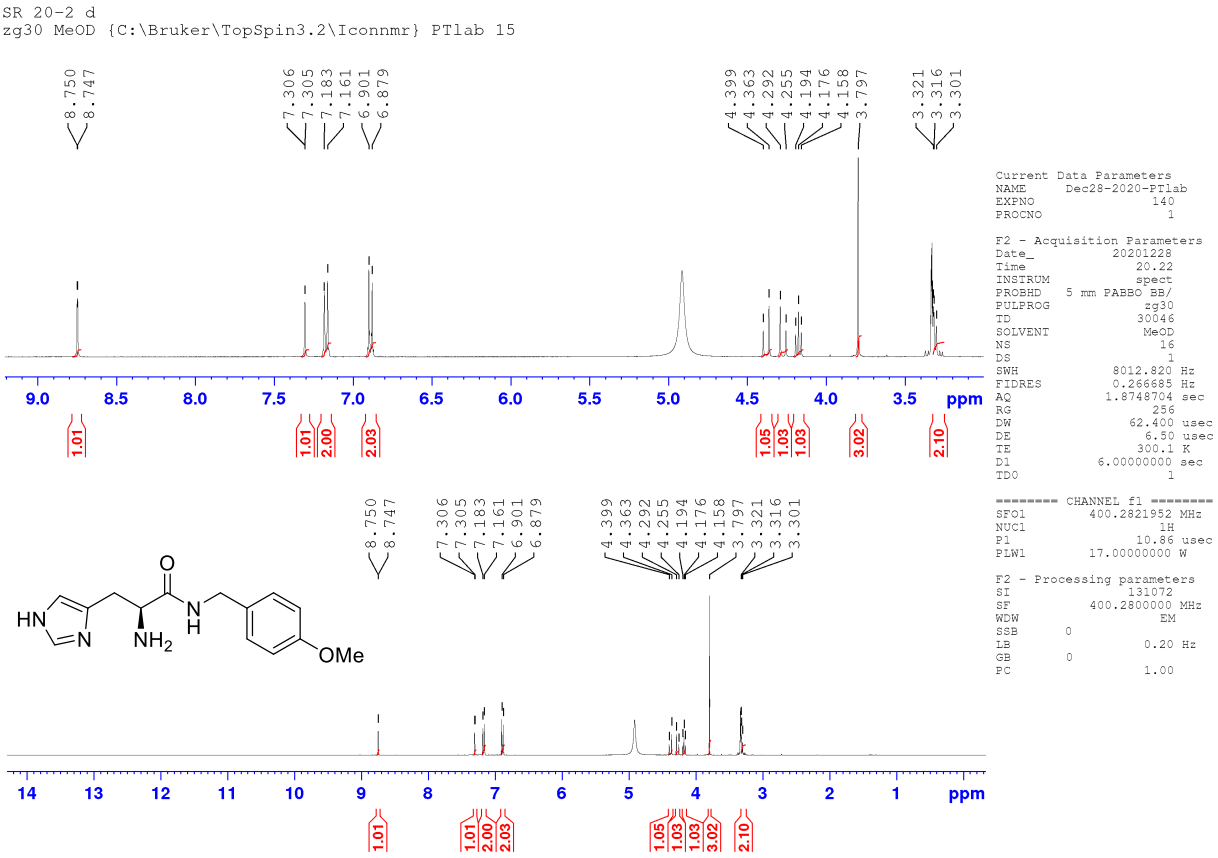
^{1}H NMR (400 MHz, MeOH-d$_4$), (*S*)-2-amino-*N*-benzyl-3-(1*H*-imidazol-4-yl)propanamide **(5a):**



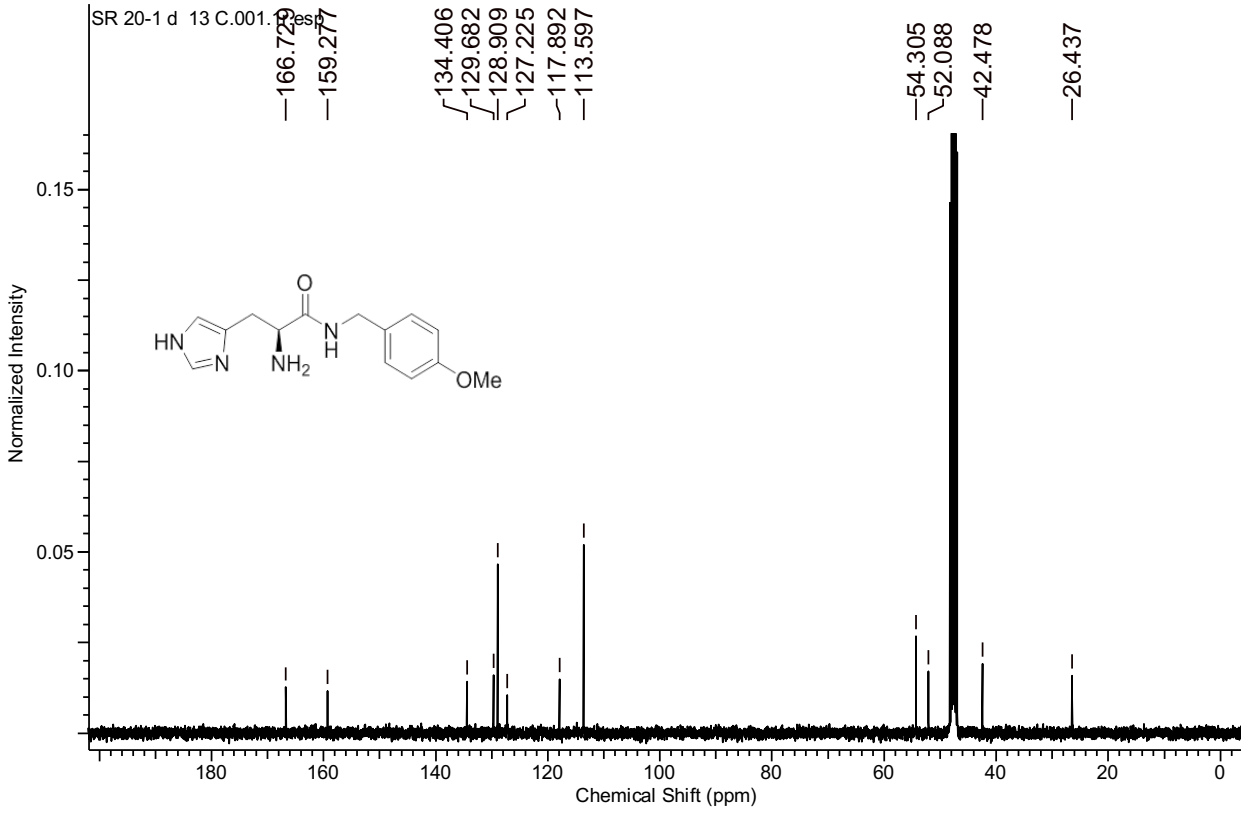
^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-*N*-benzyl-3-(1*H*-imidazol-4-yl)propanamide **(5a):**



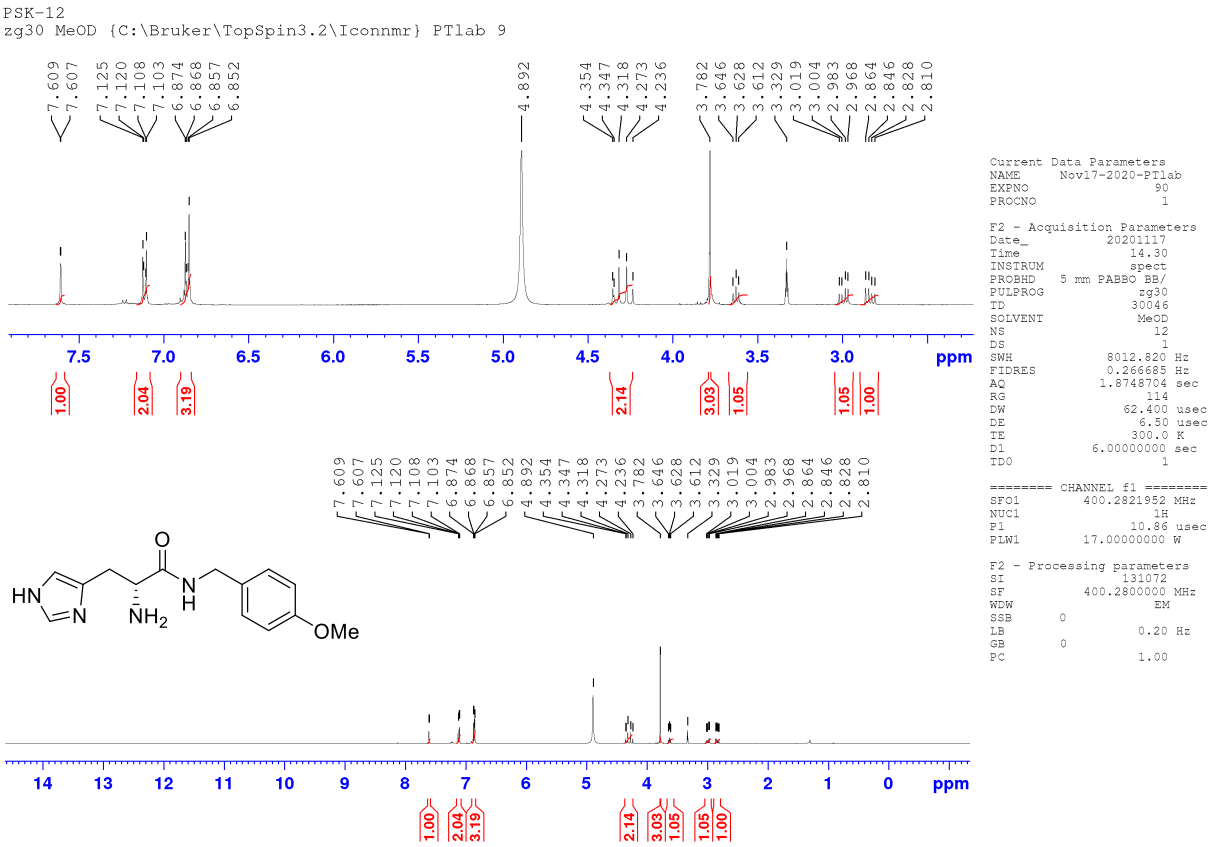
^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(4-methoxybenzyl)propanamide (**5b**):



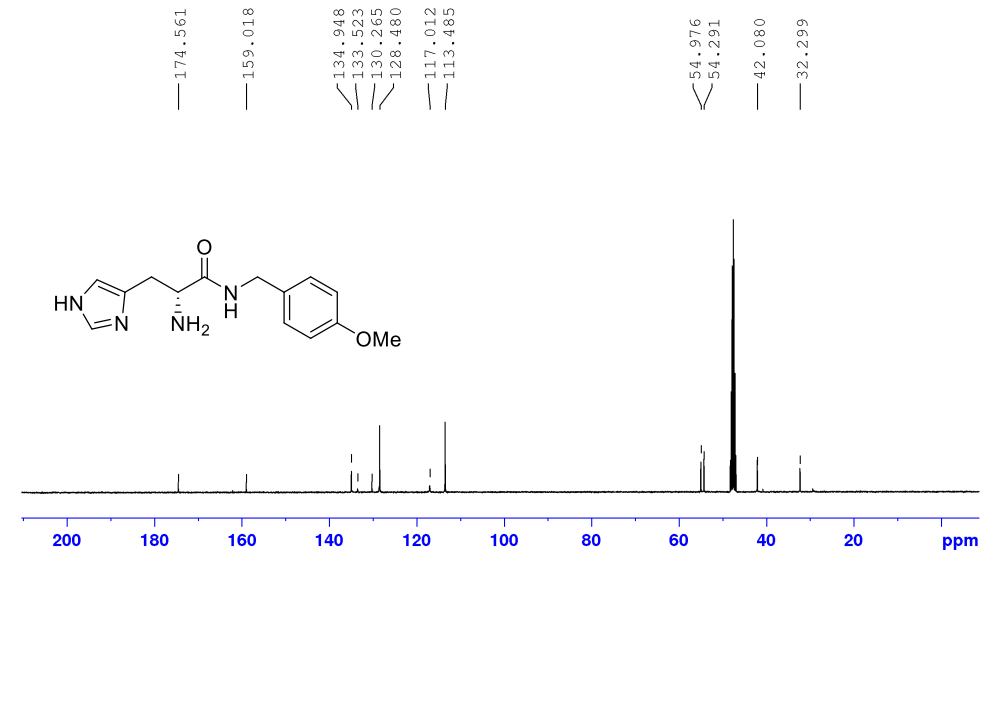
^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(4-methoxybenzyl)propanamide (**5b):**



^{1}H NMR (400 MHz, MeOH-d$_4$), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-methoxybenzyl)propanamide**(5c):**



^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-methoxybenzyl)propanamide**(5c):**




```
Current Data Parameters
NAME      Nov17-2020-PTlab
EXPNO                  150
PROCNO                   1

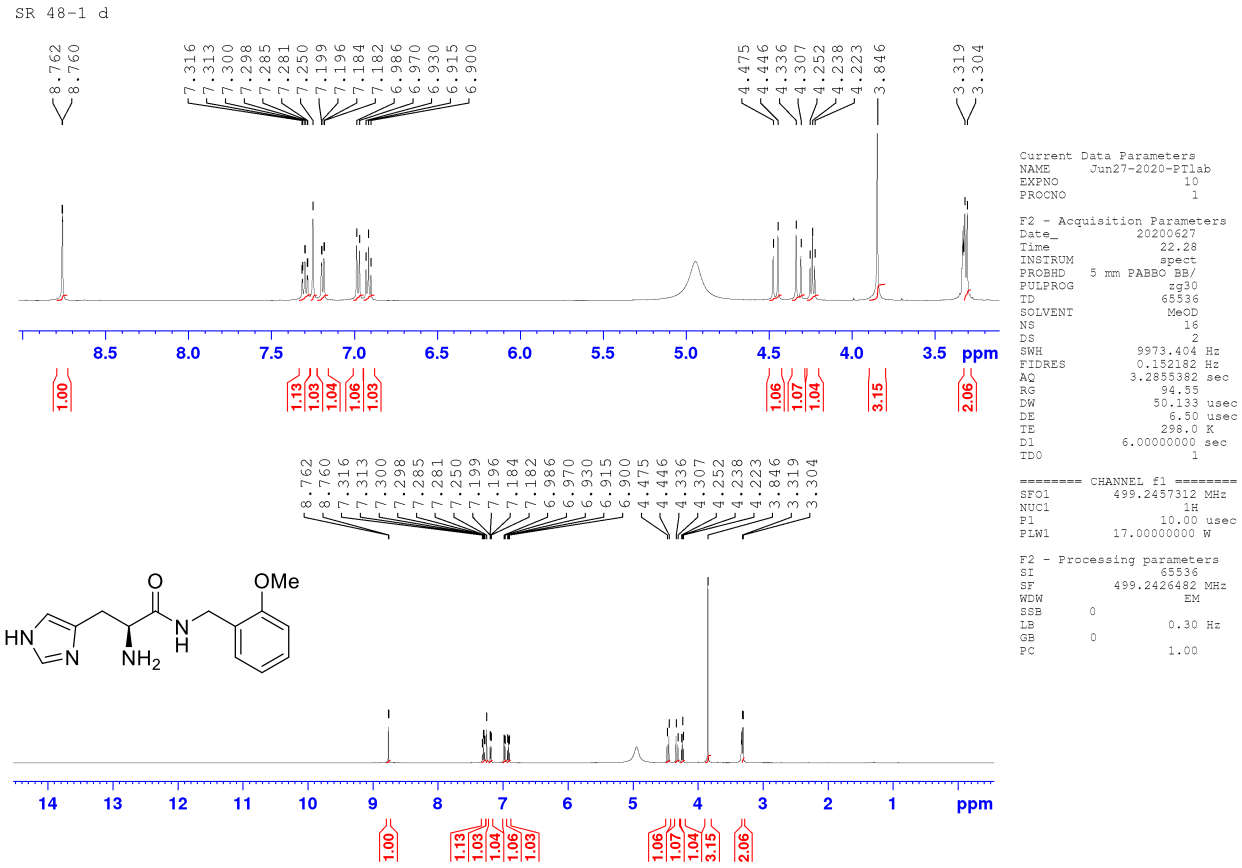
F2 - Acquisition Parameters
Date_              20201117
Time                 18.13
INSTRUM              spect
PROBHD    5 mm PABBO BB/
PULPROG            zgpg30
TD                  32768
SOLVENT              MeOD
NS                   1024
DS                      4
SWH            22058.824 Hz
FIDRES          0.673182 Hz
AQ             0.7427413 sec
RG                   1620
DW               22.667 usec
DE                 6.50 usec
TE                 300.1 K
D1            1.50000000 sec
D11           0.03000000 sec
TD0                     1

======== CHANNEL f1 ========
SFO1         100.6606368 MHz
NUC1                   13C
P1                   11.00 usec
PLW1          48.00000000 W

======== CHANNEL f2 ========
SFO2         400.2818195 MHz
NUC2                   1H
CPDPRG[2           waltz16
PCPD2                90.00 usec
PLW2          17.00000000 W
PLW12          0.24753000 W
PLW13          0.20050000 W

F2 - Processing parameters
SI                   16384
SF           100.6504861 MHz
WDW                    EM
SSB       0
LB                   1.50 Hz
GB        0
PC                   1.40
```

^{1}H NMR (500 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(2-methoxybenzyl)propanamide (**5d**):



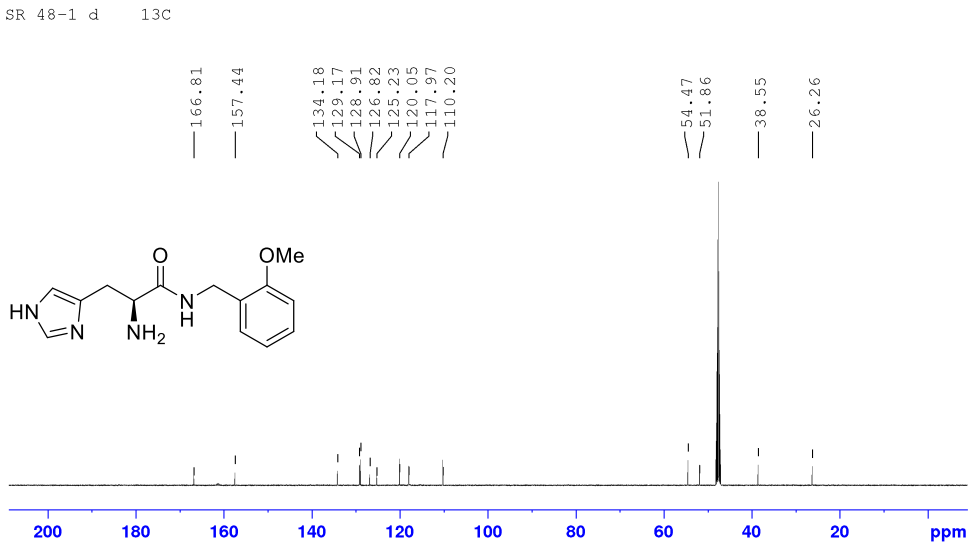

```
Current Data Parameters
NAME      Jun27-2020-PTlab
EXPNO                   10
PROCNO                   1

F2 - Acquisition Parameters
Date_            20200627
Time                22.28
INSTRUM             spect
PROBHD    5 mm PABBO BB/
PULPROG              zg30
TD                  65536
SOLVENT              MeOD
NS                     16
DS                      2
SWH              9973.404 Hz
FIDRES           0.152182 Hz
AQ               3.2855382 sec
RG                  94.55
DW                 50.133 usec
DE                  6.50 usec
TE                 298.0 K
D1              6.00000000 sec
TD0                     1

======== CHANNEL f1 ========
SFO1          499.2457312 MHz
NUC1                   1H
P1                  10.00 usec
PLW1          17.00000000 W

F2 - Processing parameters
SI                  65536
SF            499.2426482 MHz
WDW                    EM
SSB       0
LB                  0.30 Hz
GB        0
PC                  1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(2-methoxybenzyl)propanamide (**5d**):




```
Current Data Parameters
NAME      Jun27-2020-PTlab
EXPNO                   20
PROCNO                   1

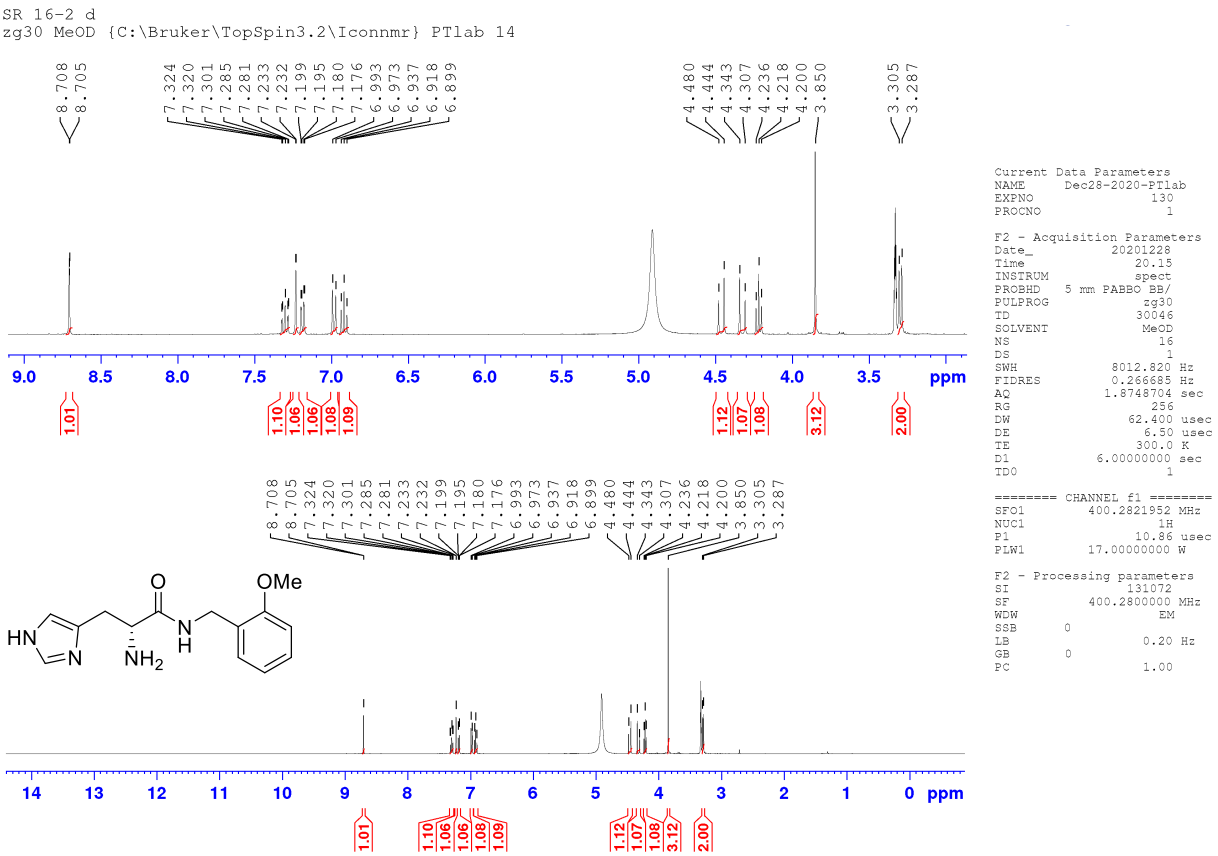
F2 - Acquisition Parameters
Date_               20200627
Time                   23.03
INSTRUM                spect
PROBHD   5 mm PABBO BB/
PULPROG             zgpg30
TD                   65536
SOLVENT                MeOD
NS                       512
DS                         4
SWH              29761.904 Hz
FIDRES            0.454131 Hz
AQ                1.1010048 sec
RG                   191.52
DW                 16.800 usec
DE                 10.00 usec
TE                   300.0 K
D1              2.00000000 sec
D11             0.03000000 sec
TD0                      1

======== CHANNEL f1 ========
SFO1          125.5472166 MHz
NUC1                    13C
P1                  10.00 usec
PLW1          71.00000000 W

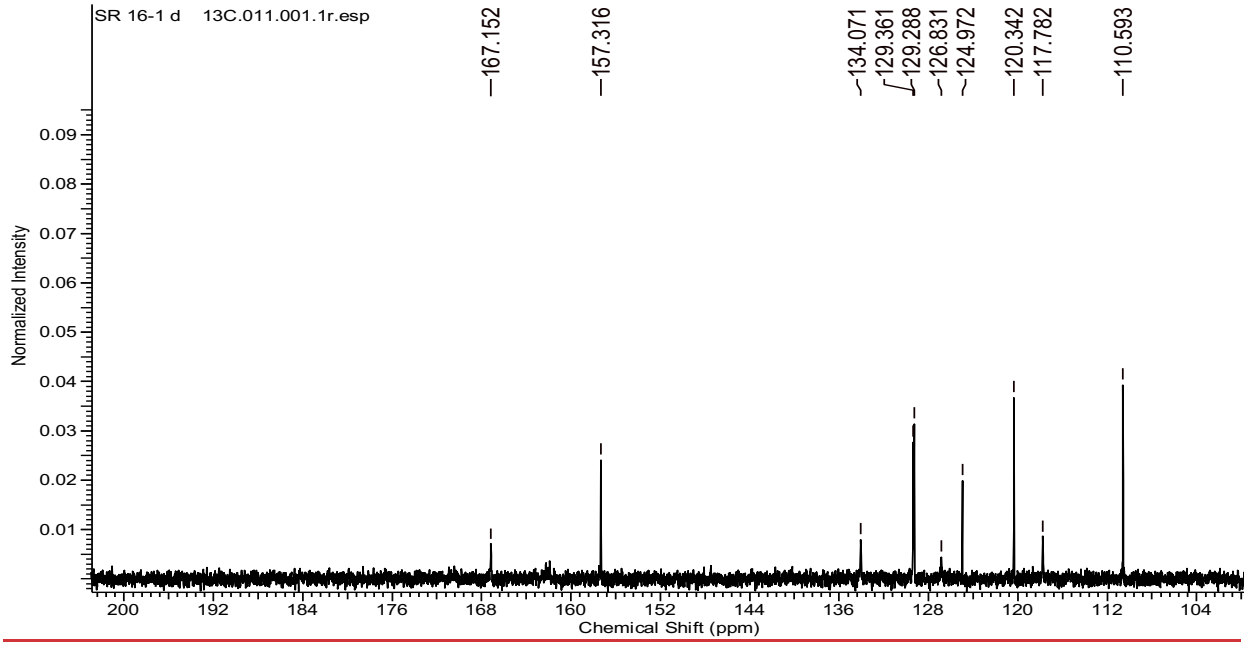
======== CHANNEL f2 ========
SFO2          499.2446452 MHz
NUC2                    1H
CPDPRG[2           waltz16
PCPD2               80.00 usec
PLW2          17.00000000 W
PLW12          0.26563001 W
PLW13          0.17000000 W

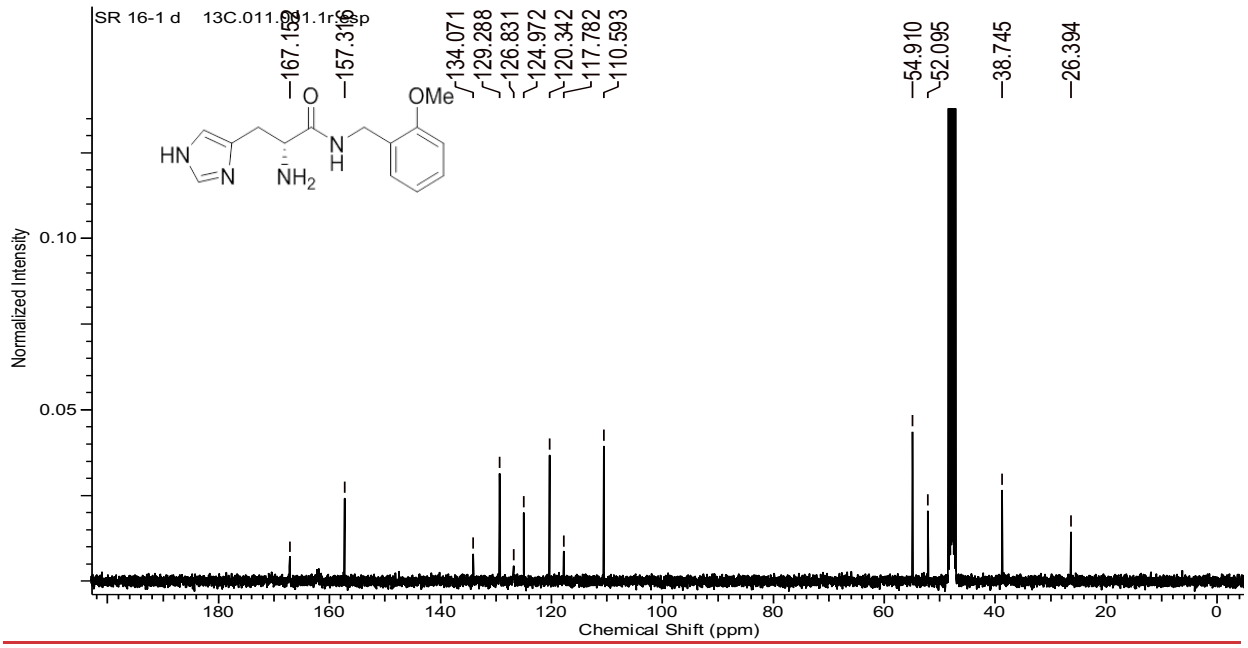
F2 - Processing parameters
SI                     32768
SF            125.5346637 MHz
WDW                       EM
SSB       0
LB                     1.00 Hz
GB        0
PC                     1.40
```

^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(2-methoxybenzyl)propanamide (**5e):**

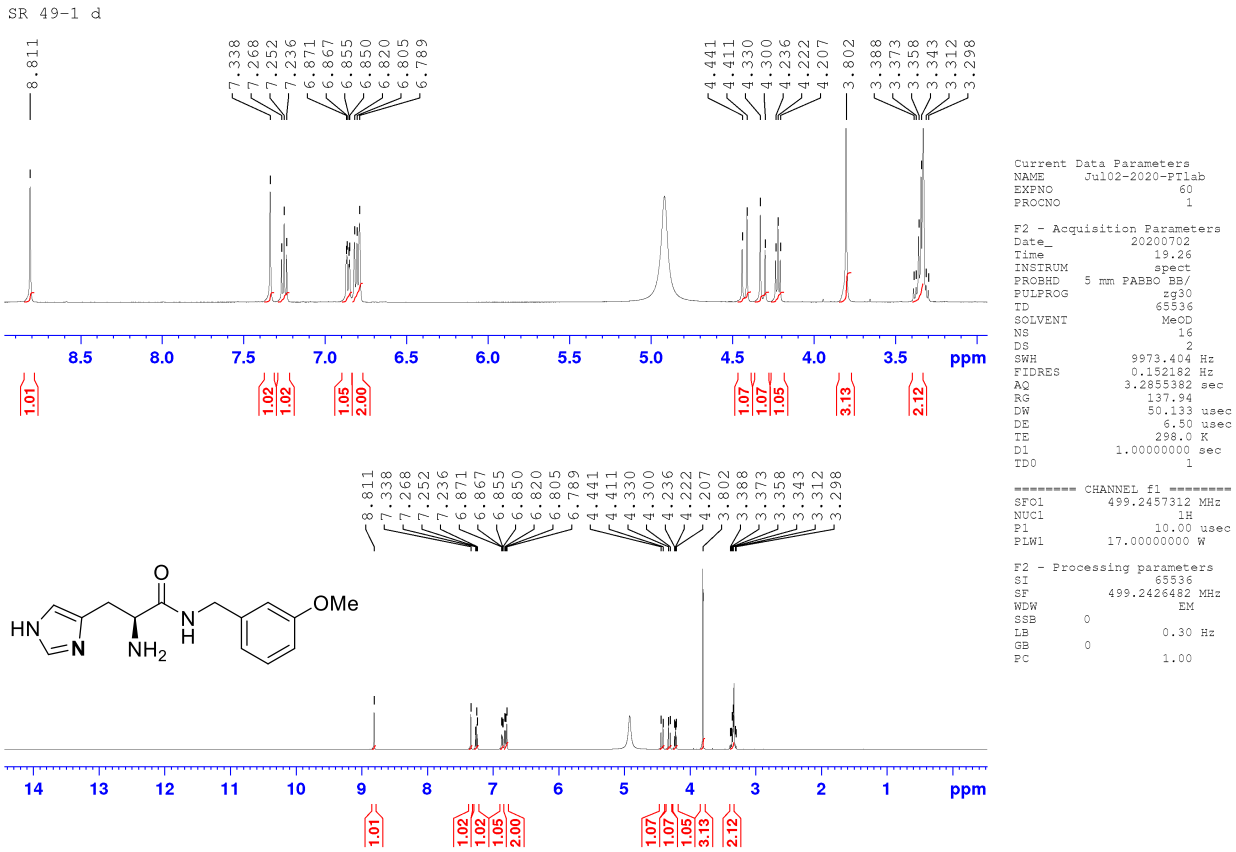


^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(2-methoxybenzyl)propanamide (**5e):**

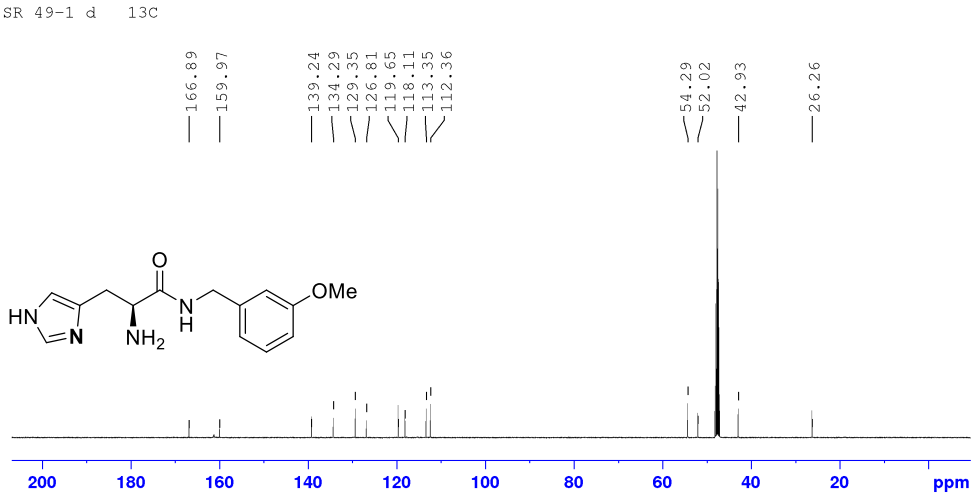




^{1}H NMR (500 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(3-methoxybenzyl)propanamide (**5f**)**:**



^{13}C NMR (126 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(3-methoxybenzyl)propanamide (**5f**)**:**




```
Current Data Parameters
NAME       Jul02-2020-PTlab
EXPNO                   70
PROCNO                   1

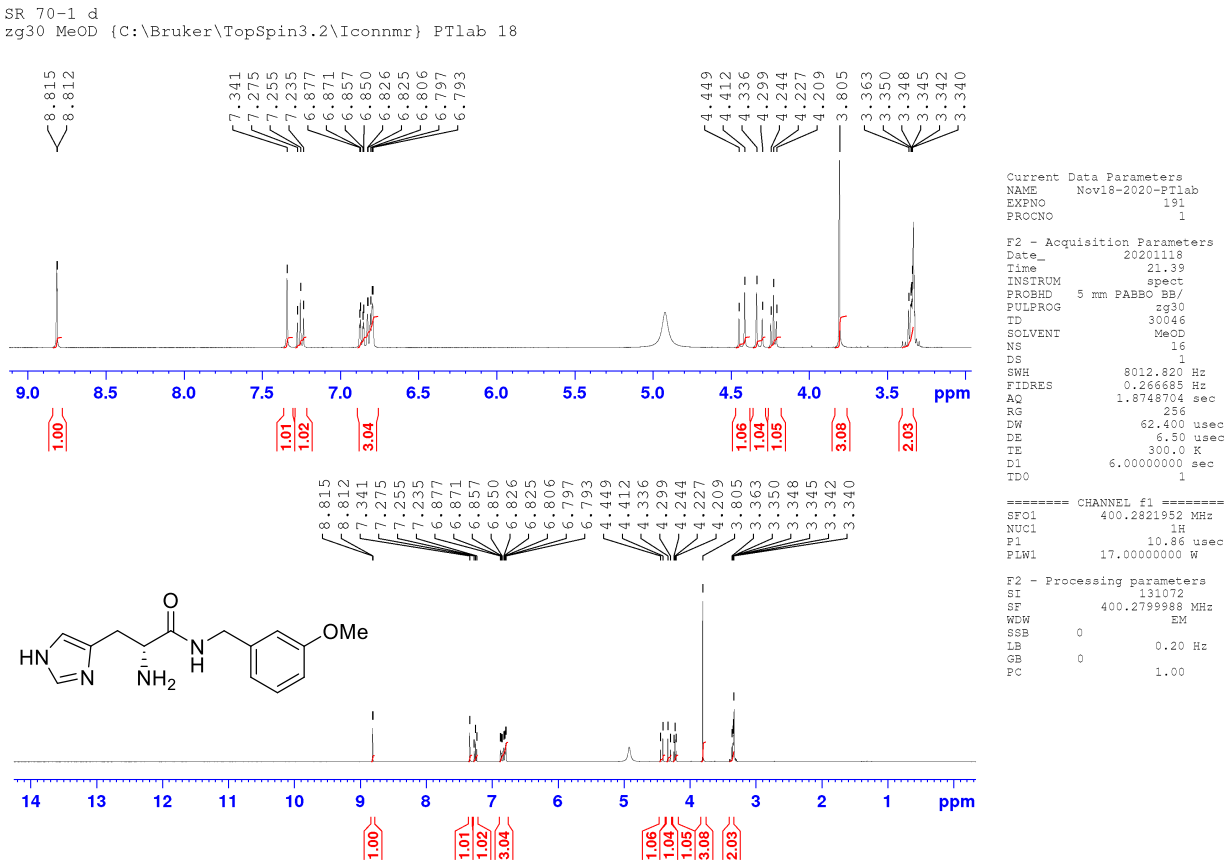
F2 - Acquisition Parameters
Date_              20200702
Time                 20.11
INSTRUM              spect
PROBHD    5 mm PABBO BB/
PULPROG           zgpg30
TD                 65536
SOLVENT              MeOD
NS                     512
DS                       4
SWH            29761.904 Hz
FIDRES          0.454131 Hz
AQ             1.1010048 sec
RG                  191.52
DW                16.800 usec
DE                10.00 usec
TE                300.0 K
D1             2.00000000 sec
D11            0.03000000 sec
TD0                      1

======== CHANNEL f1 ========
SFO1         125.5472166 MHz
NUC1                   13C
P1                   10.00 usec
PLW1         71.00000000 W

======== CHANNEL f2 ========
SFO2         499.2446452 MHz
NUC2                   1H
CPDPRG[2          waltz16
PCPD2                80.00 usec
PLW2         17.00000000 W
PLW12         0.26563001 W
PLW13         0.17000000 W

F2 - Processing parameters
SI                   32768
SF           125.5346637 MHz
WDW                     EM
SSB       0
LB                   1.00 Hz
GB        0
PC                   1.40
```

^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(3-methoxybenzyl)propanamide (**5g**):



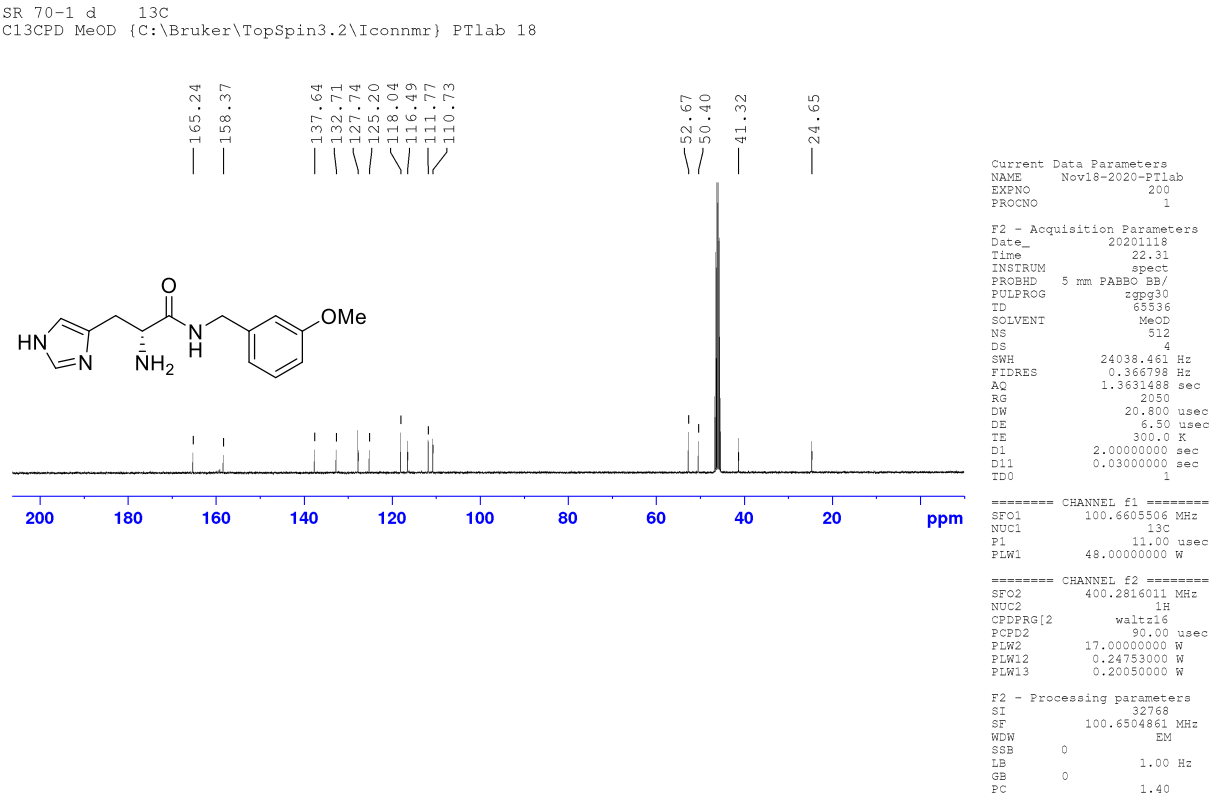

```
Current Data Parameters
NAME      Nov18-2020-PTlab
EXPNO                  191
PROCNO                   1

F2 - Acquisition Parameters
Date_           20201118
Time              21.39
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG            zg30
TD                30046
SOLVENT            MeOD
NS                   16
DS                    1
SWH            8012.820 Hz
FIDRES         0.266685 Hz
AQ             1.8748704 sec
RG                  256
DW                62.400 usec
DE                 6.50 usec
TE                300.0 K
D1             6.00000000 sec
TD0                   1

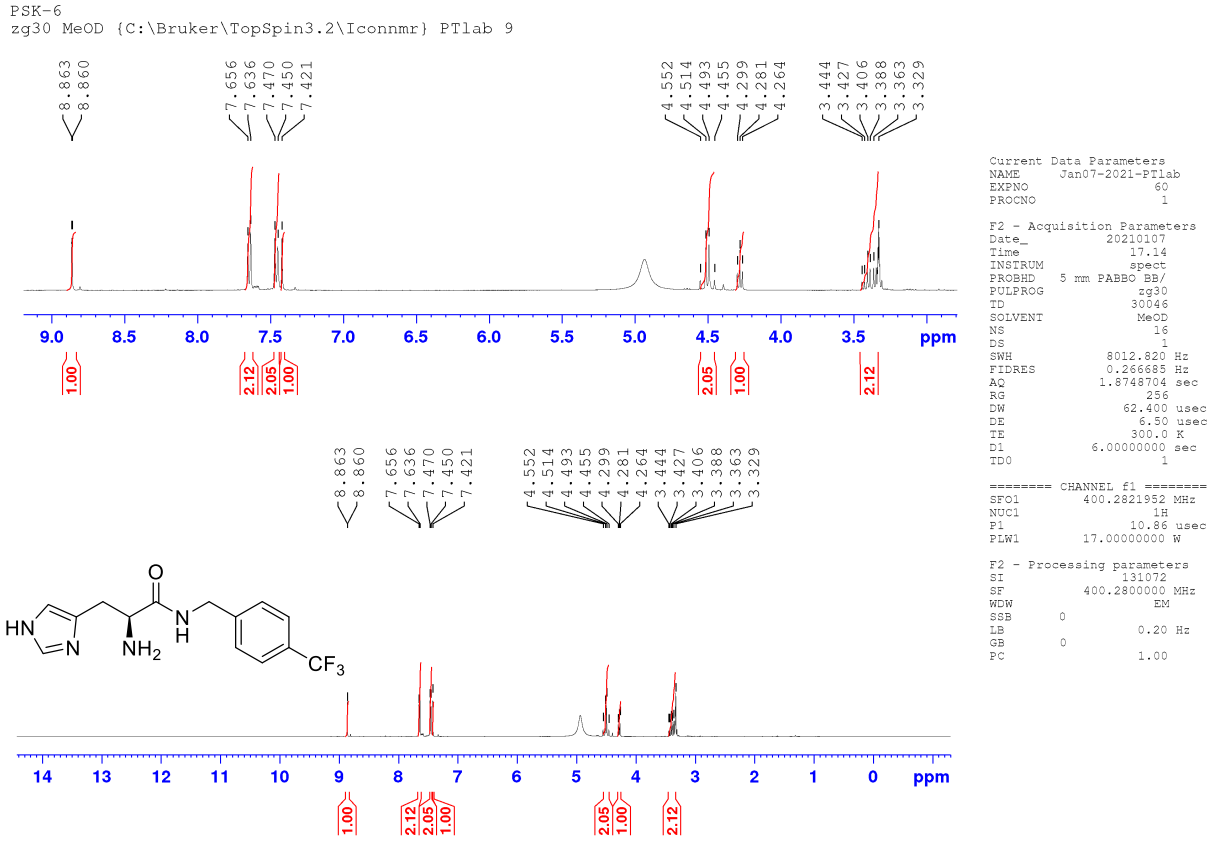
======== CHANNEL f1 ========
SFO1         400.2821952 MHz
NUC1                    1H
P1                  10.86 usec
PLW1        17.00000000 W

F2 - Processing parameters
SI                  131072
SF           400.2799988 MHz
WDW                    EM
SSB       0
LB                  0.20 Hz
GB        0
PC                  1.00
```

^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(3-methoxybenzyl)propanamide (**5g**):



^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)benzyl)propanamide **(5h):**



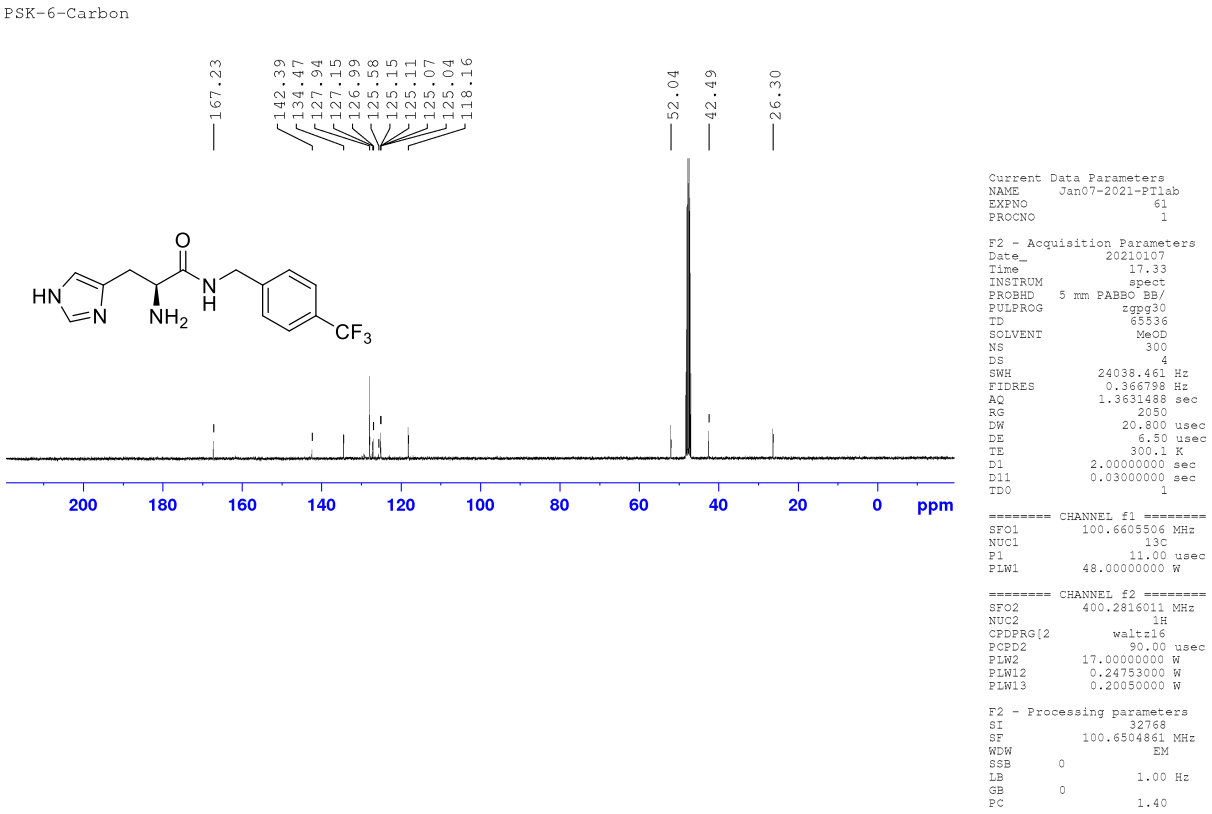

```
Current Data Parameters
NAME      Jan07-2021-PTlab
EXPNO                   60
PROCNO                   1

F2 - Acquisition Parameters
Date_           20210107
Time              17.14
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG           zg30
TD               30046
SOLVENT           MeOD
NS                  16
DS                   1
SWH            8012.820 Hz
FIDRES         0.266685 Hz
AQ             1.8748704 sec
RG                 256
DW              62.400 usec
DE               6.50 usec
TE              300.0 K
D1            6.00000000 sec
TD0                  1

======== CHANNEL f1 ========
SFO1         400.2821952 MHz
NUC1               1H
P1              10.86 usec
PLW1        17.00000000 W

F2 - Processing parameters
SI              131072
SF           400.2800000 MHz
WDW                EM
SSB       0
LB              0.20 Hz
GB        0
PC              1.00
```

^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)benzyl)propanamide **(5h):**




```
Current Data Parameters
NAME      Jan07-2021-PTlab
EXPNO                   61
PROCNO                   1

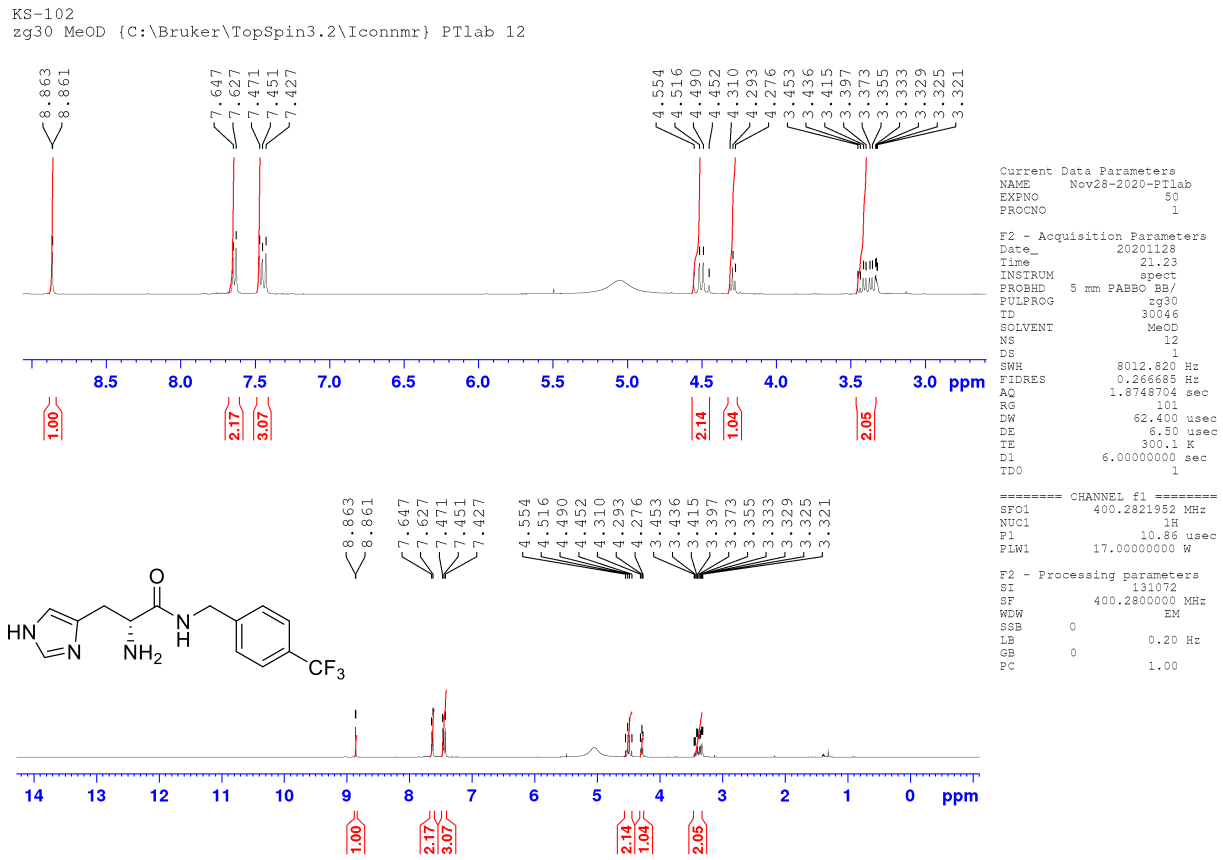
F2 - Acquisition Parameters
Date_           20210107
Time              17.33
INSTRUM           spect
PROBHD    5 mm PABBO BB/
PULPROG          zgpg30
TD                65536
SOLVENT            MeOD
NS                  300
DS                    4
SWH           24038.461 Hz
FIDRES         0.366798 Hz
AQ             1.3631488 sec
RG                 2050
DW                20.800 usec
DE                 6.50 usec
TE                300.1 K
D1            2.00000000 sec
D11           0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1         100.6605506 MHz
NUC1                 13C
P1                 11.00 usec
PLW1         48.00000000 W

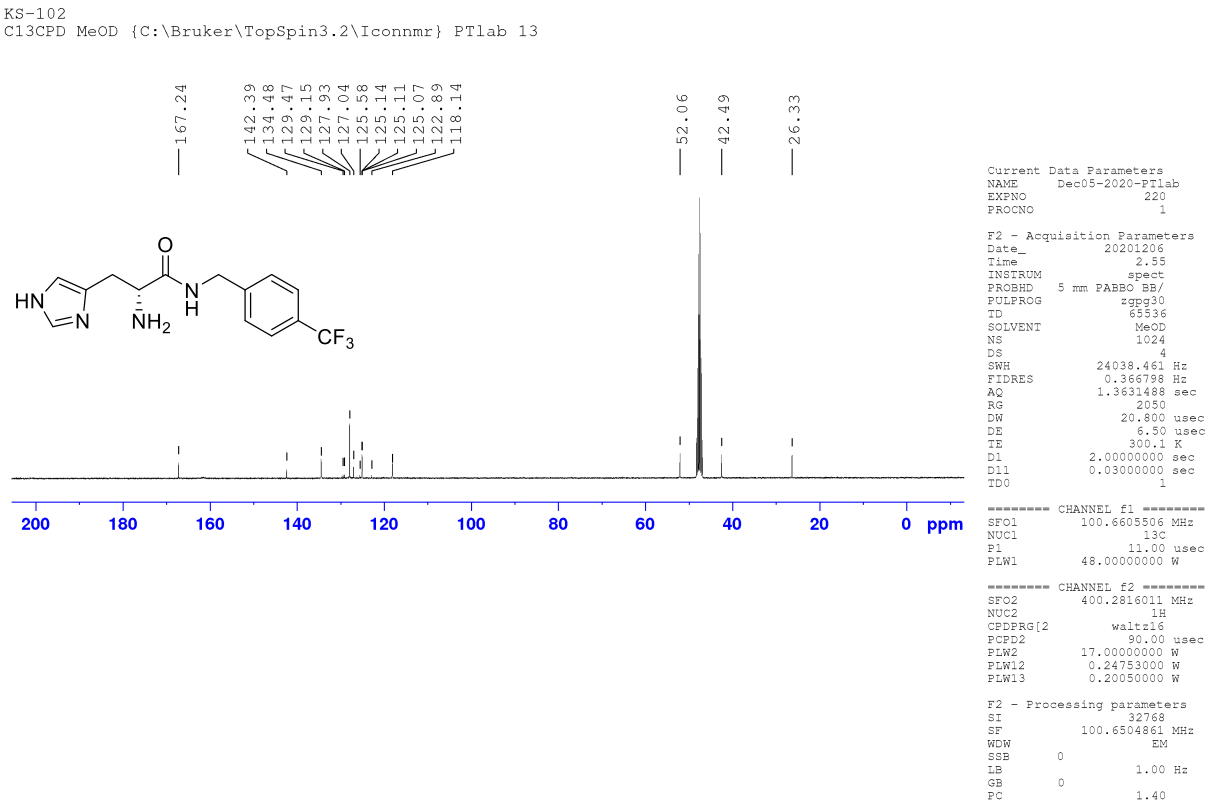
======== CHANNEL f2 ========
SFO2         400.2816011 MHz
NUC2                 1H
CPDPRG[2          waltz16
PCPD2              90.00 usec
PLW2         17.00000000 W
PLW12         0.24753000 W
PLW13         0.20050000 W

F2 - Processing parameters
SI                  32768
SF           100.6504861 MHz
WDW                  EM
SSB       0
LB                 1.00 Hz
GB        0
PC                 1.40
```

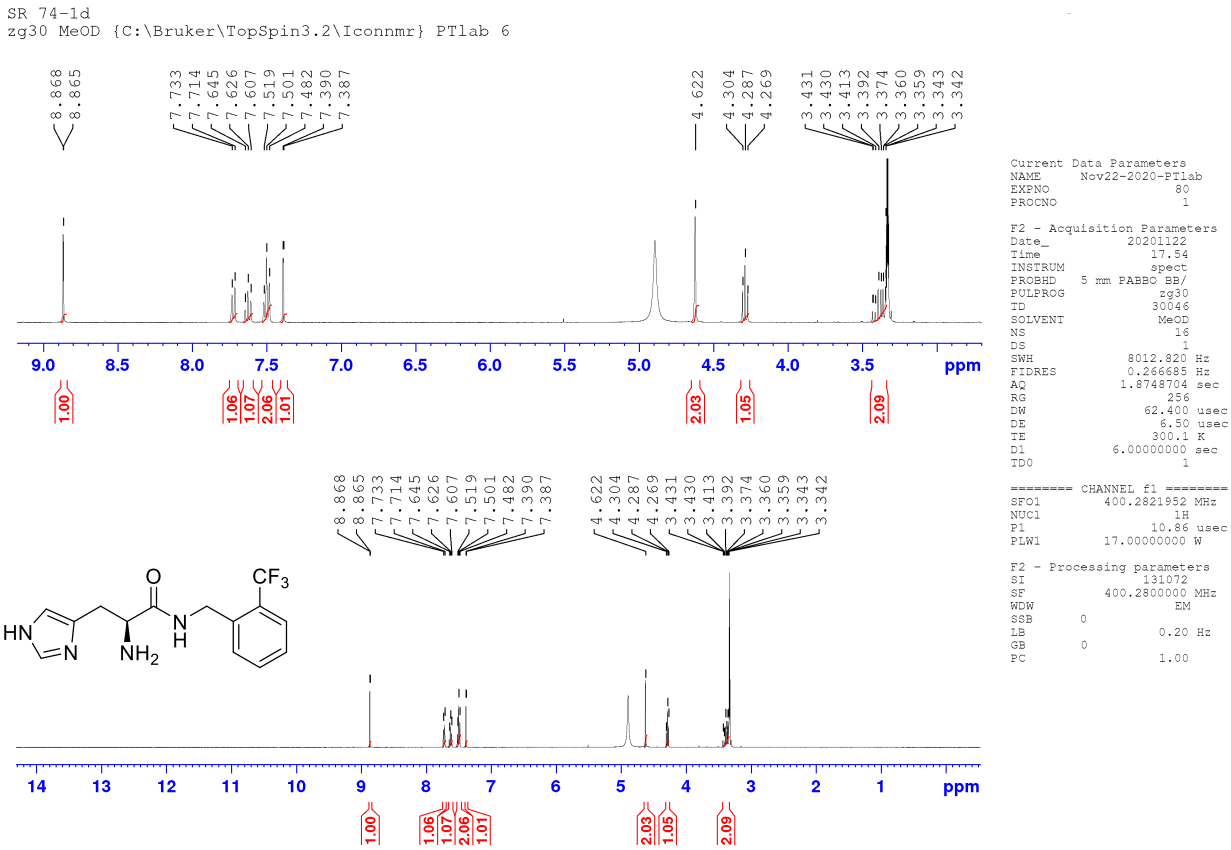
^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)benzyl)propanamide **(5i):**



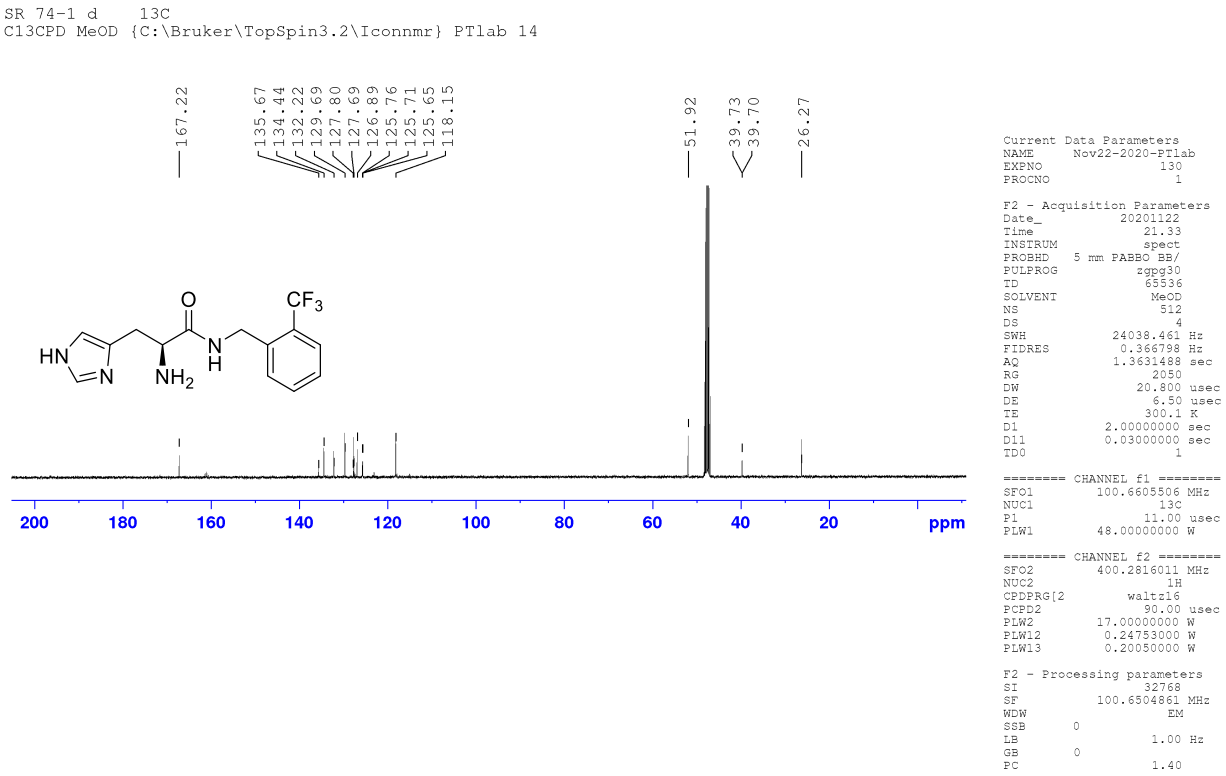
^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)benzyl)propanamide **(5i):**



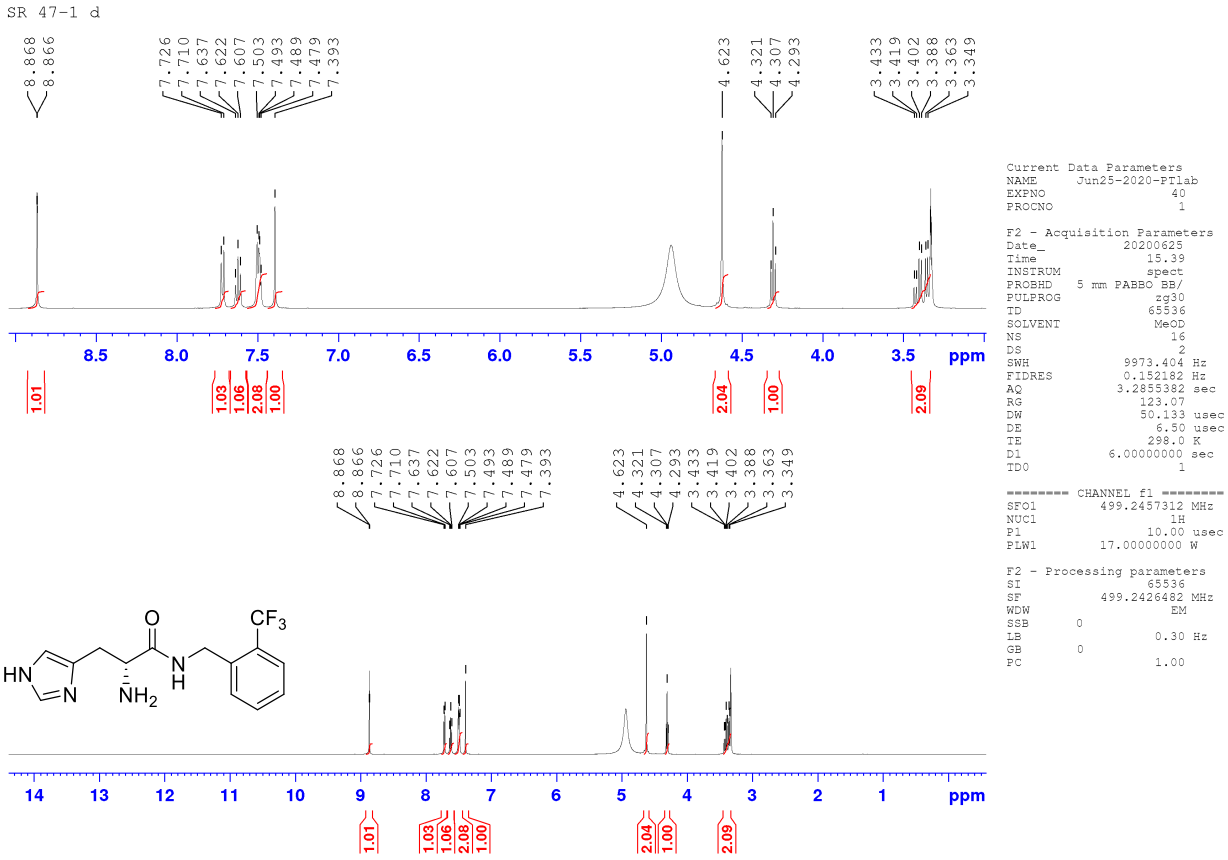
^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(trifluoromethyl)benzyl)propanamide (**5j**):



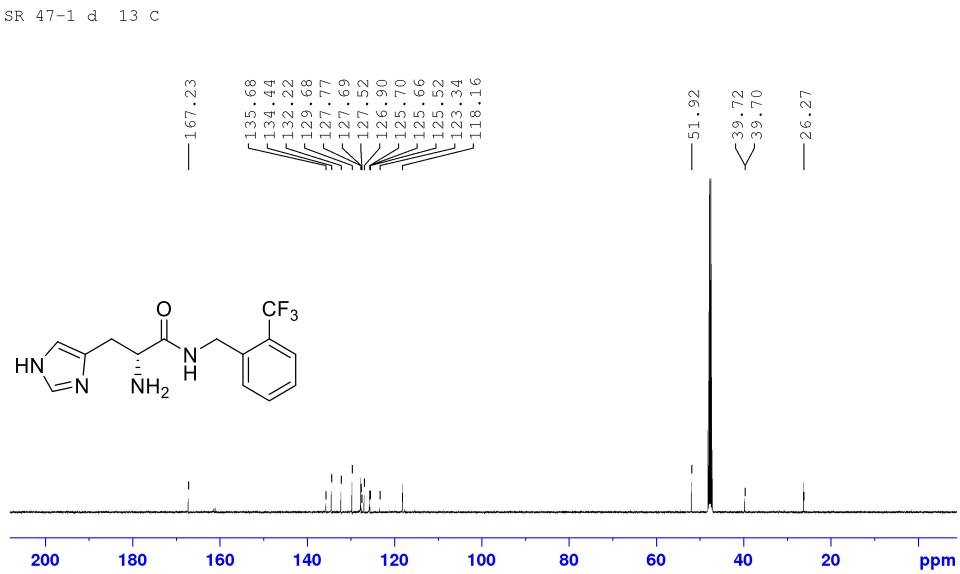
^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(trifluoromethyl)benzyl)propanamide (**5j**):



^{1}H NMR (500 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(trifluoromethyl)benzyl)propanamide (**5k**):



^{13}C NMR (126 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(trifluoromethyl)benzyl)propanamide (**5k**):




```
Current Data Parameters
NAME      Jun25-2020-PTlab
EXPNO                   50
PROCNO                   1

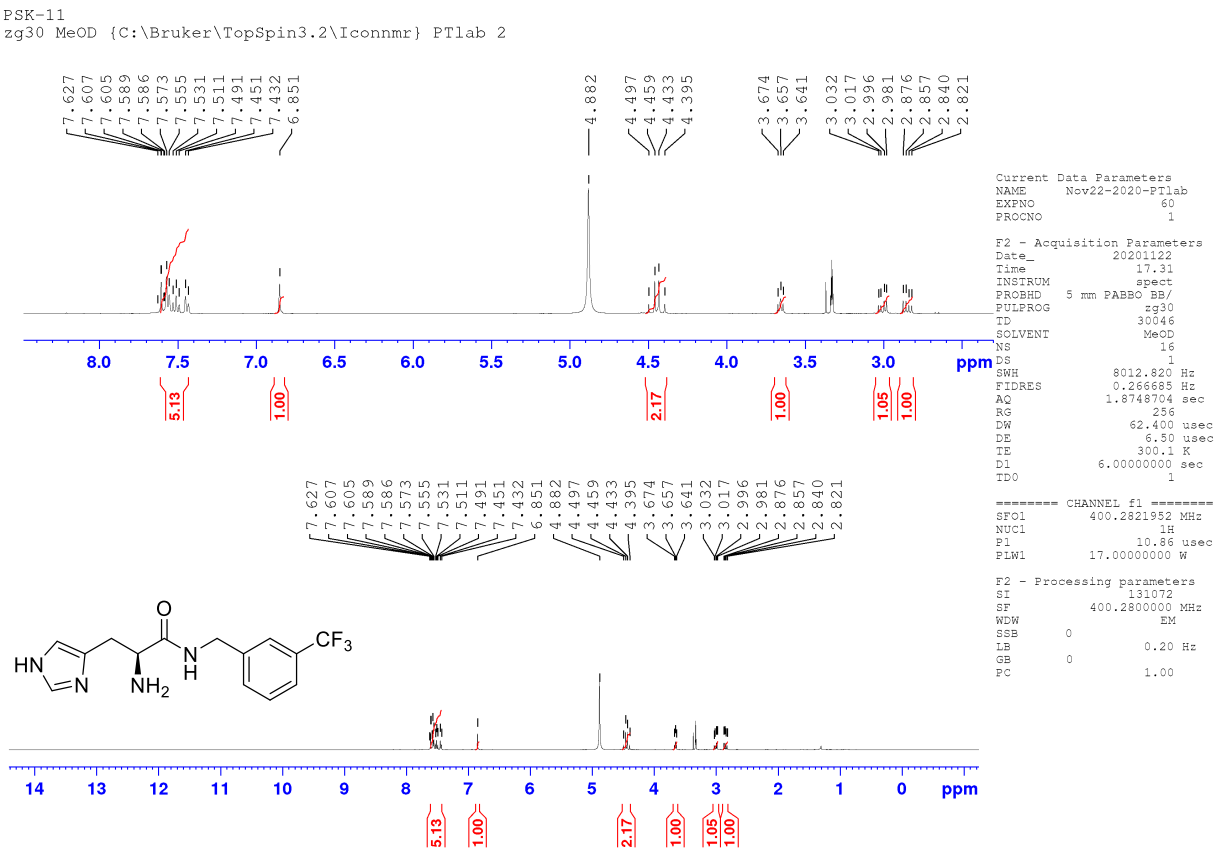
F2 - Acquisition Parameters
Date_           20200625
Time              16.02
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG         zgpg30
TD               65536
SOLVENT            MeOD
NS                  450
DS                    4
SWH           29761.904 Hz
FIDRES         0.454131 Hz
AQ             1.1010048 sec
RG                191.52
DW                16.800 usec
DE                10.00 usec
TE                300.0 K
D1            2.00000000 sec
D11           0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1         125.5472166 MHz
NUC1                13C
P1                10.00 usec
PLW1         71.00000000 W

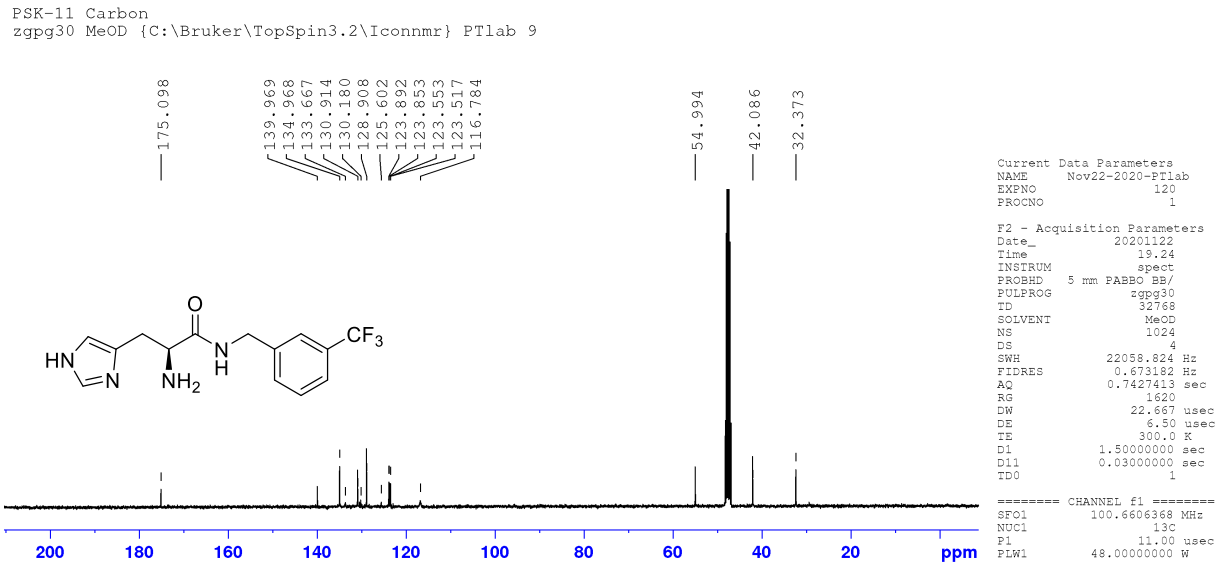
======== CHANNEL f2 ========
SFO2         499.2446452 MHz
NUC2                1H
CPDPRG[2         waltz16
PCPD2             80.00 usec
PLW2         17.00000000 W
PLW12         0.26563001 W
PLW13         0.17000000 W

F2 - Processing parameters
SI                32768
SF           125.5346637 MHz
WDW                  EM
SSB       0
LB                1.00 Hz
GB        0
PC                1.40
```

^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-(trifluoromethyl)benzyl)propanamide **(5l):**



^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-(trifluoromethyl)benzyl)propanamide **(5l):**




```
Current Data Parameters
NAME       Nov22-2020-PT1ab
EXPNO                  120
PROCNO                   1

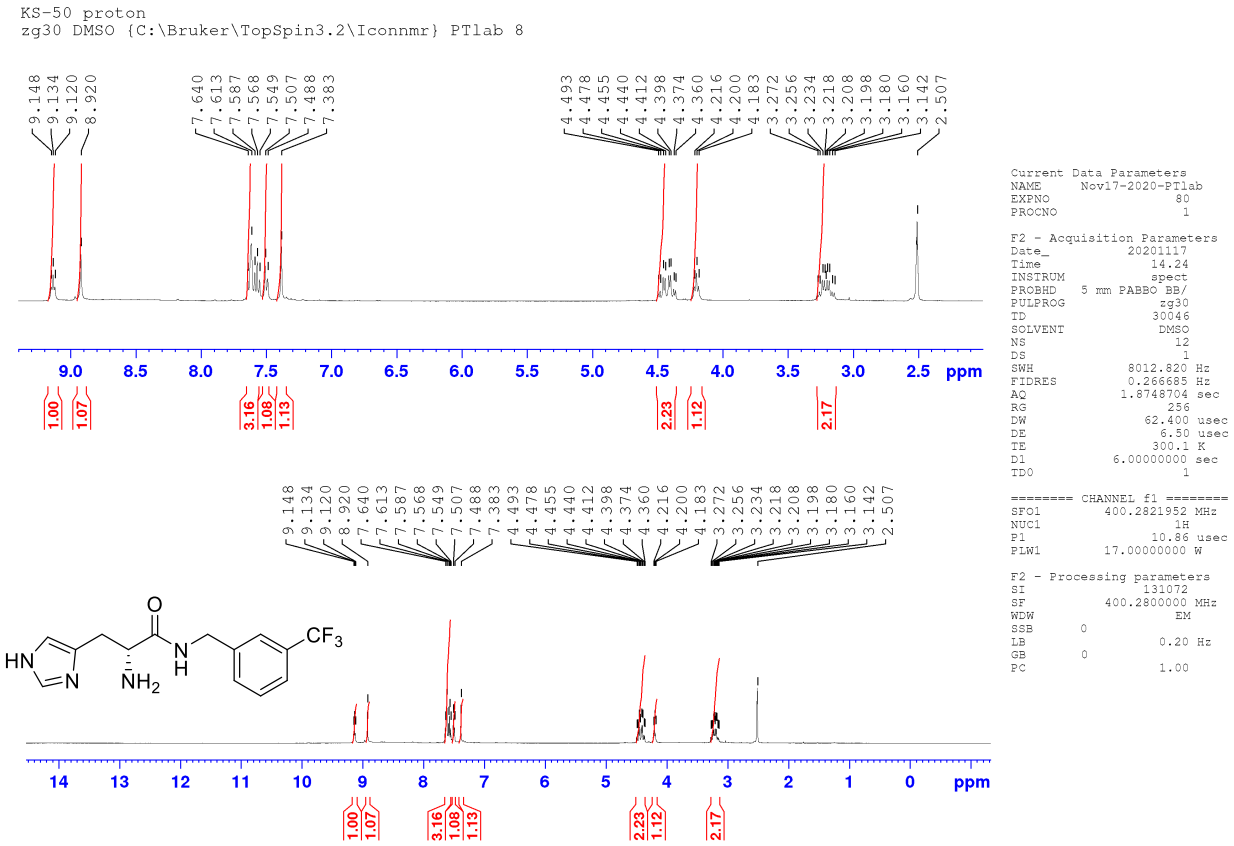
F2 - Acquisition Parameters
Date_           20201122
Time              19.24
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG          zgpg30
TD                32768
SOLVENT             MeOD
NS                  1024
DS                     4
SWH           22058.824 Hz
FIDRES         0.673182 Hz
AQ            0.7427413 sec
RG                  1620
DW               22.667 usec
DE                6.50 usec
TE                300.0 K
D1            1.50000000 sec
D11           0.03000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1         100.6606368 MHz
NUC1                13C
P1                11.00 usec
PLW1         48.00000000 W

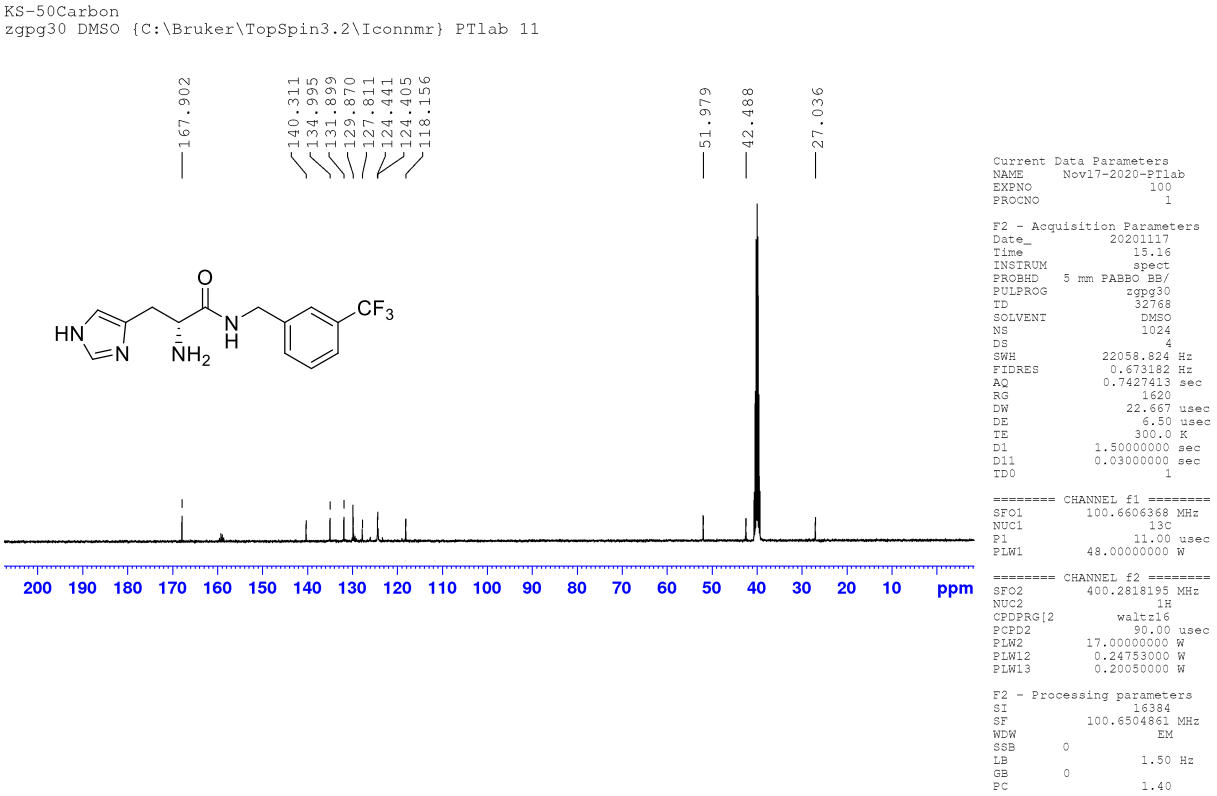
======== CHANNEL f2 ========
SFO2         400.2818195 MHz
NUC2                 1H
CPDPRG[2         waltz16
PCPD2              90.00 usec
PLW2         17.00000000 W
PLW12         0.24753000 W
PLW13         0.20050000 W

F2 - Processing parameters
SI                 16384
SF           100.6504861 MHz
WDW                   EM
SSB       0
LB                 1.50 Hz
GB        0
PC                 1.40
```

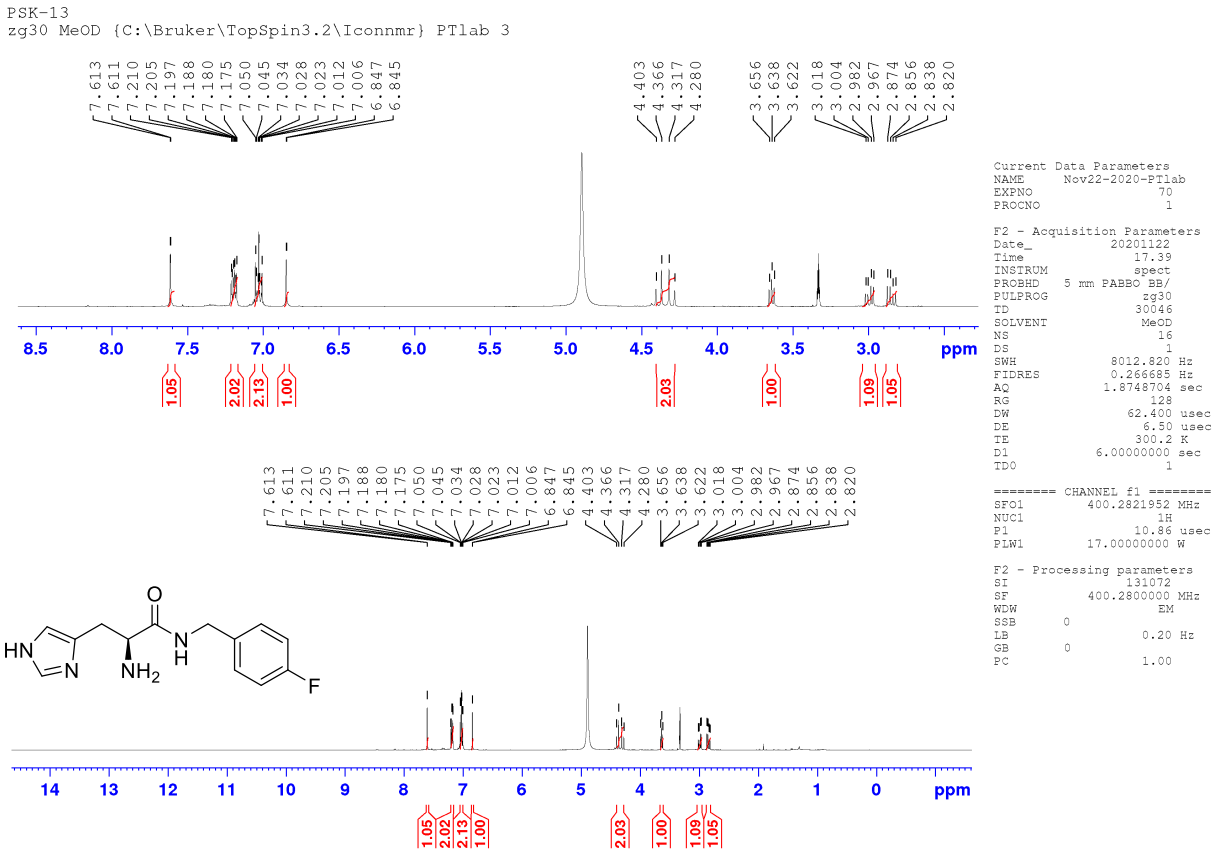
^{1}H NMR (400 MHz, DMSO-d_6), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-(trifluoromethyl)benzyl)propanamide **(5m):**



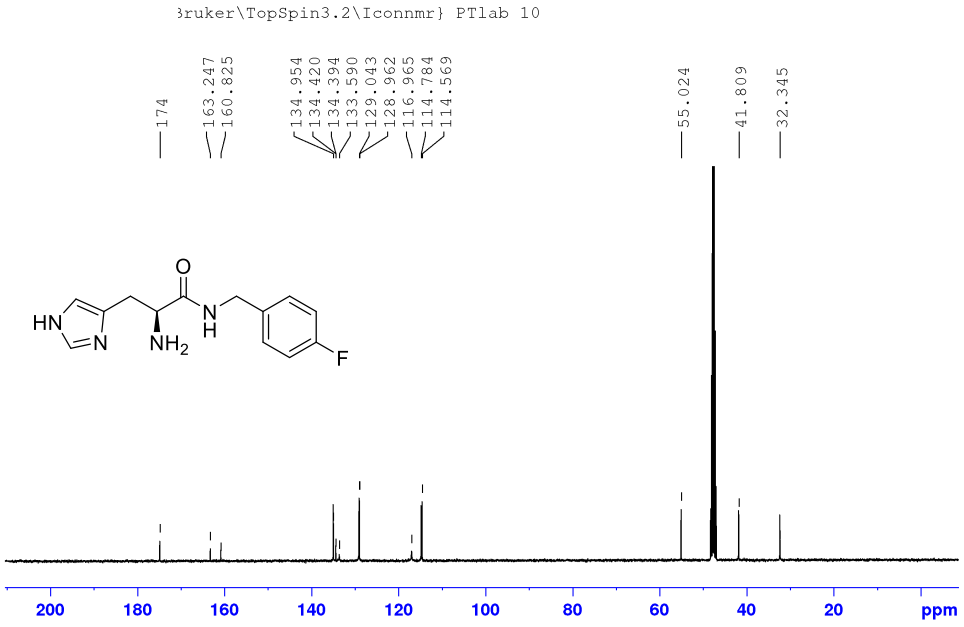
^{13}C NMR (100 MHz, DMSO-d_6), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-(trifluoromethyl)benzyl)propanamide **(5m):**



^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-*N*-(4-fluorobenzyl)-3-(1*H*-imidazol-4-yl)propanamide **(5n):**



^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-*N*-(4-fluorobenzyl)-3-(1*H*-imidazol-4-yl)propanamide **(5n):**




```
Current Data Parameters
NAME      Nov22-2020-PTlab
EXPNO                  110
PROCNO                   1

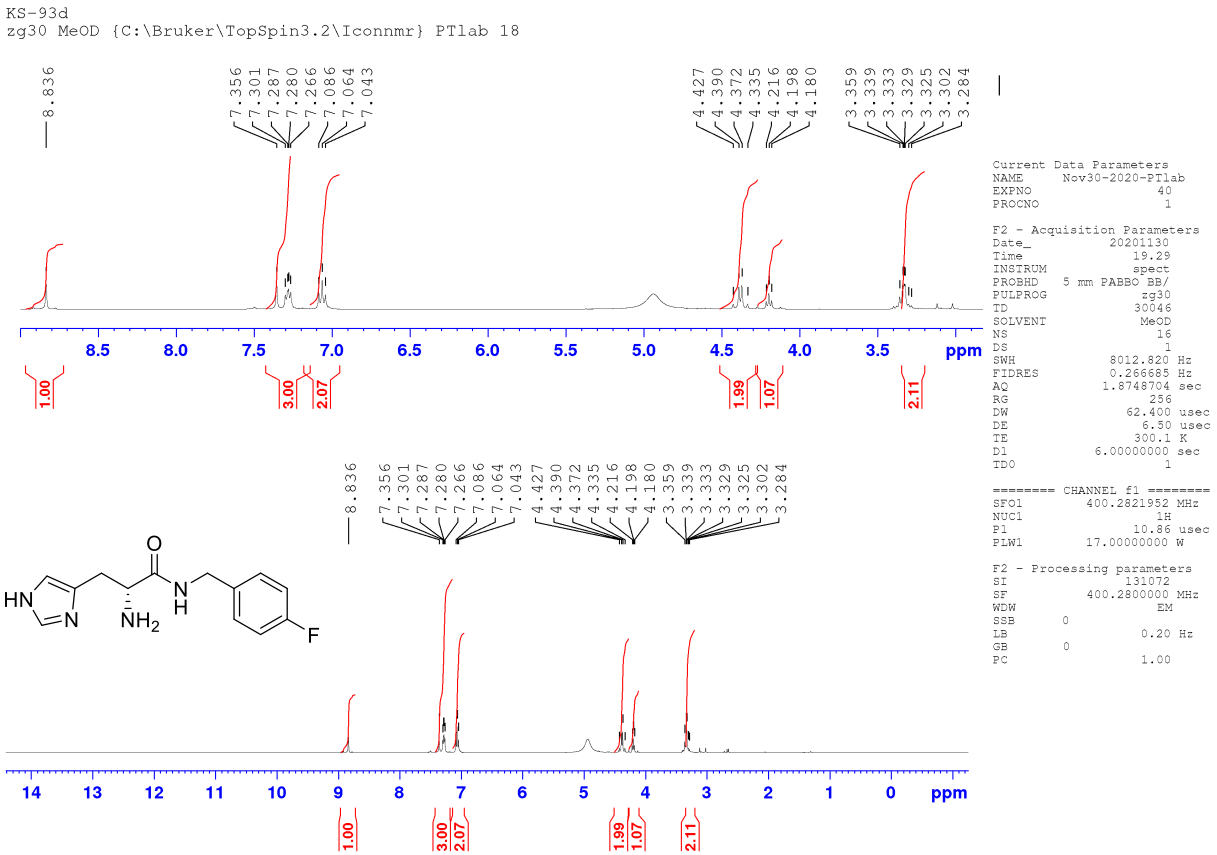
F2 - Acquisition Parameters
Date_           20201122
Time              20.09
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG         zgpg30
TD                32768
SOLVENT            MeOD
NS                 1024
DS                    4
SWH           22058.824 Hz
FIDRES         0.673182 Hz
AQ            0.7427413 sec
RG                 1620
DW               22.667 usec
DE                6.50 usec
TE                300.1 K
D1           1.50000000 sec
D11          0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1         100.6606368 MHz
NUC1                 13C
P1                11.00 usec
PLW1        48.00000000 W

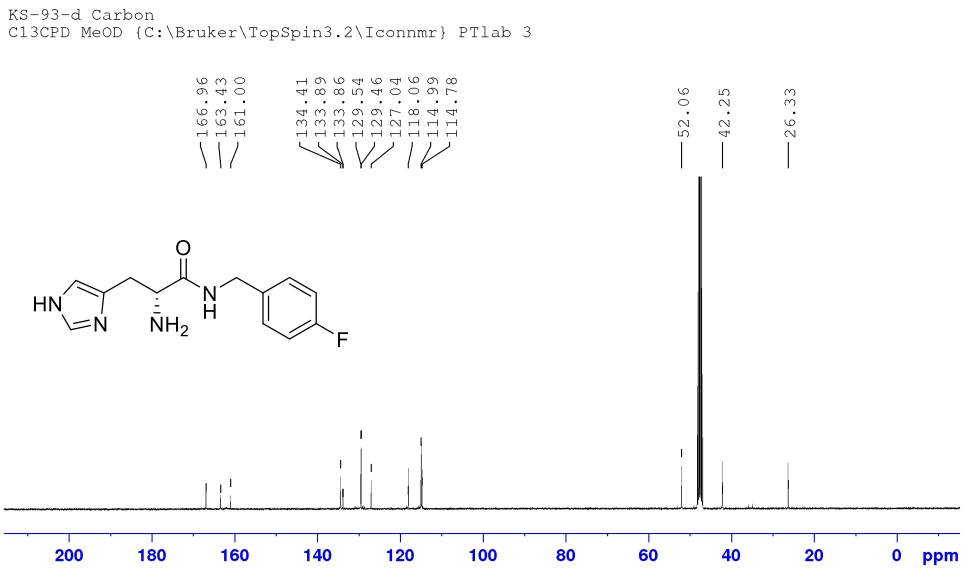
======== CHANNEL f2 ========
SFO2         400.2818195 MHz
NUC2                 1H
CPDPRG[2          waltz16
PCPD2             90.00 usec
PLW2        17.00000000 W
PLW12        0.24753000 W
PLW13        0.20050000 W

F2 - Processing parameters
SI                16384
SF           100.6504861 MHz
WDW                  EM
SSB       0
LB                1.50 Hz
GB        0
PC                1.40
```

^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-*N*-(4-fluorobenzyl)-3-(1*H*-imidazol-4-yl)propanamide **(5o):**



^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-*N*-(4-fluorobenzyl)-3-(1*H*-imidazol-4-yl)propanamide **(5o):**




```
Current Data Parameters
NAME      Jan30-2021-PTlab
EXPNO                   70
PROCNO                   1

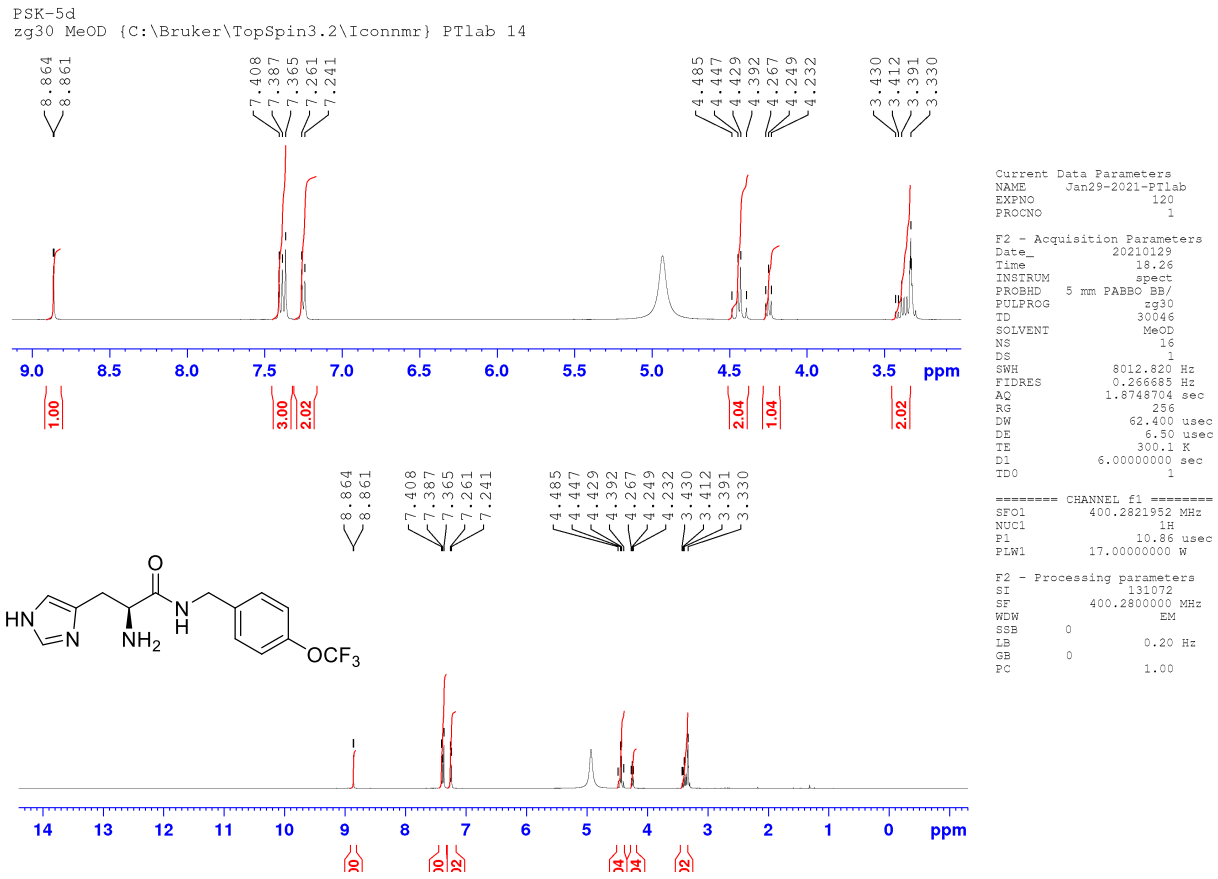
F2 - Acquisition Parameters
Date_            20210130
Time                15.48
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG          zgpg30
TD                65536
SOLVENT            MeOD
NS                 1024
DS                    4
SWH           24038.461 Hz
FIDRES         0.366798 Hz
AQ             1.3631488 sec
RG                 2050
DW               20.800 usec
DE                6.50 usec
TE                300.1 K
D1            2.00000000 sec
D11           0.03000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1         100.6605506 MHz
NUC1               13C
P1                11.00 usec
PLW1         48.00000000 W

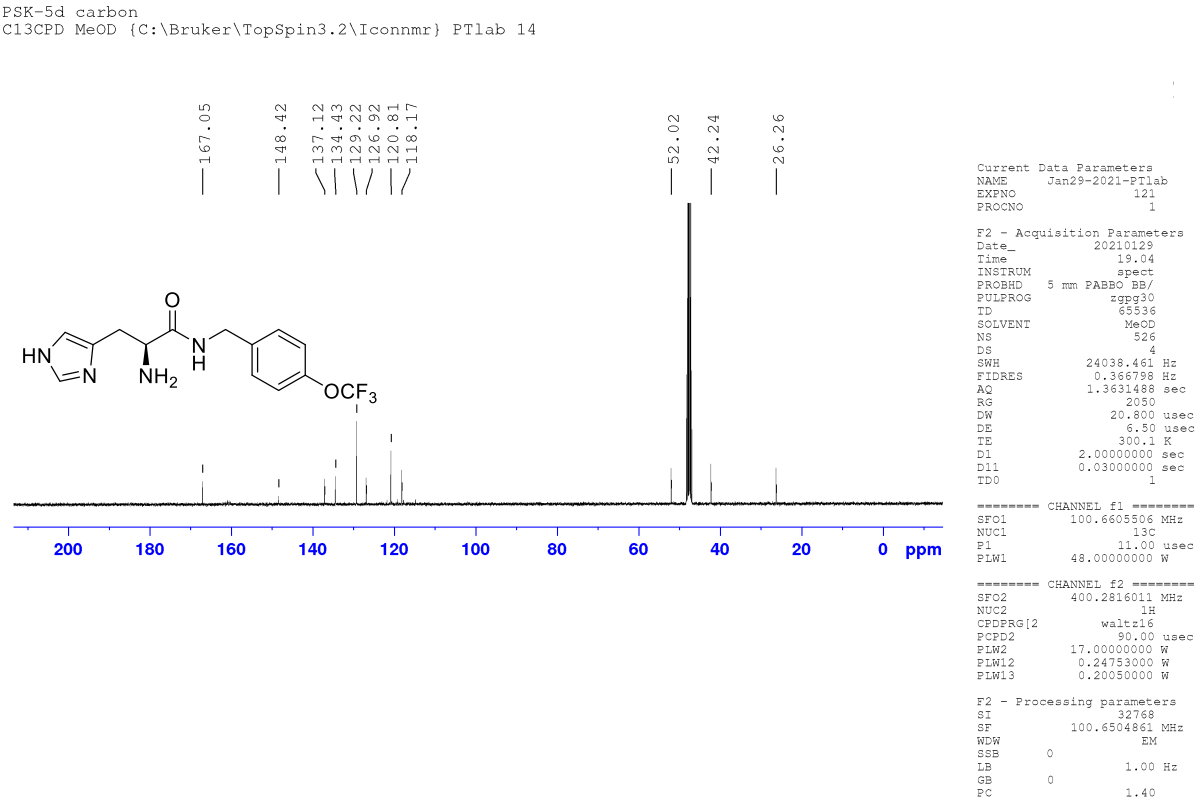
======== CHANNEL f2 ========
SFO2         400.2816011 MHz
NUC2               1H
CPDPRG[2         waltz16
PCPD2             90.00 usec
PLW2         17.00000000 W
PLW12         0.24753000 W
PLW13         0.20050000 W

F2 - Processing parameters
SI                 32768
SF           100.6504861 MHz
WDW                  EM
SSB       0
LB                 1.00 Hz
GB        0
PC                 1.40
```

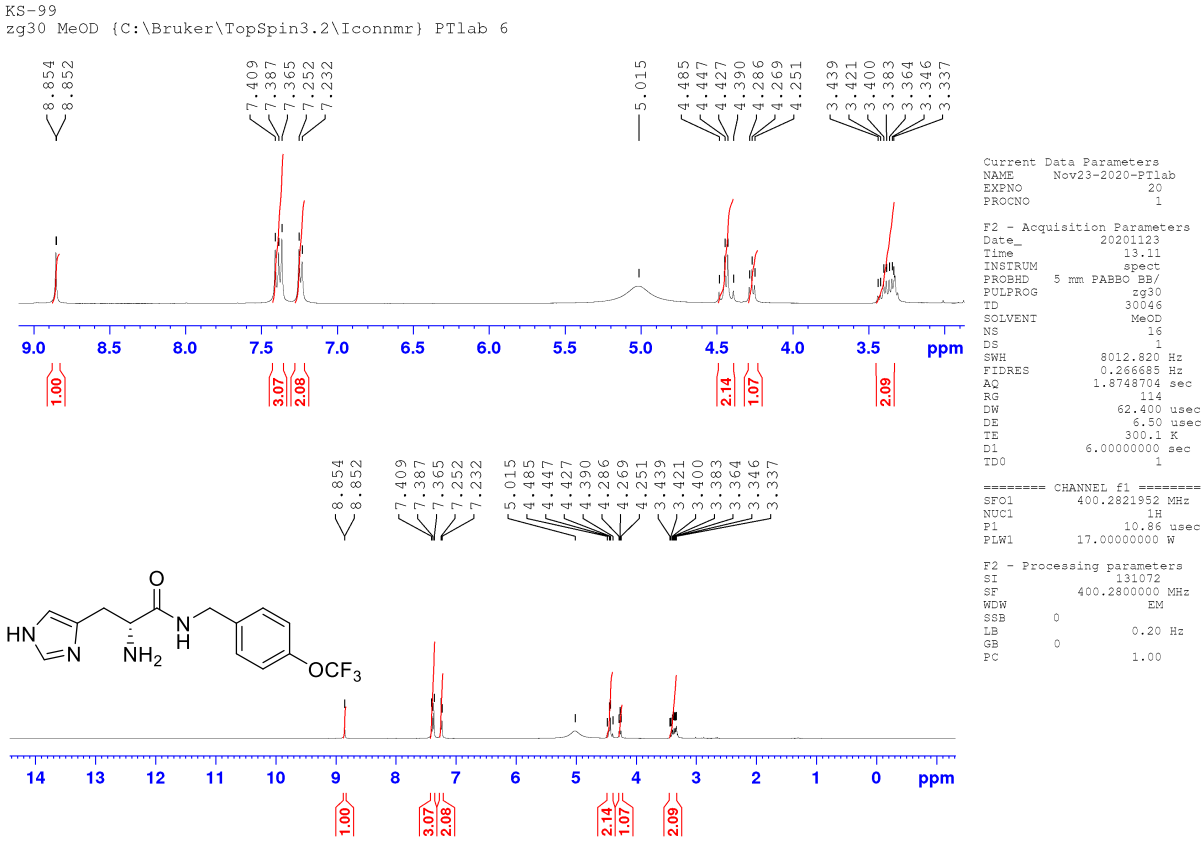
^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethoxy)benzyl)propanamide **(5p):**



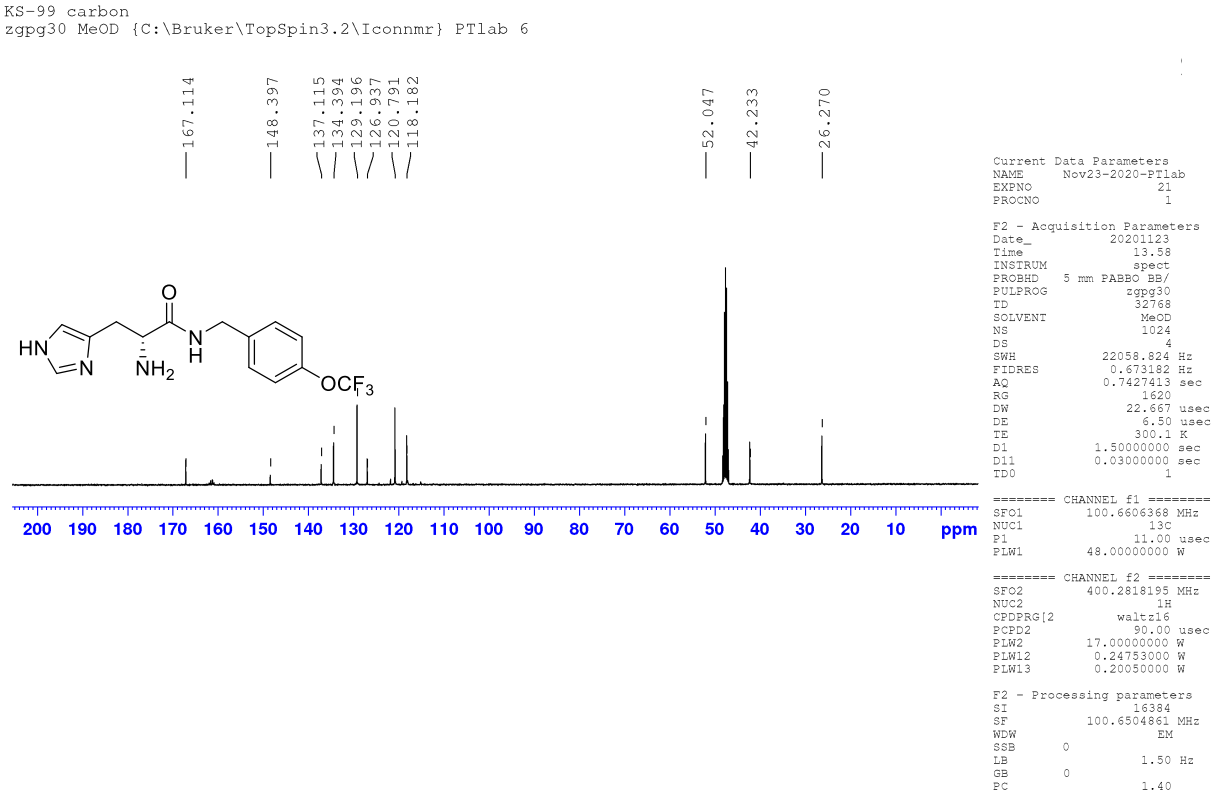
^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethoxy)benzyl)propanamide **(5p):**



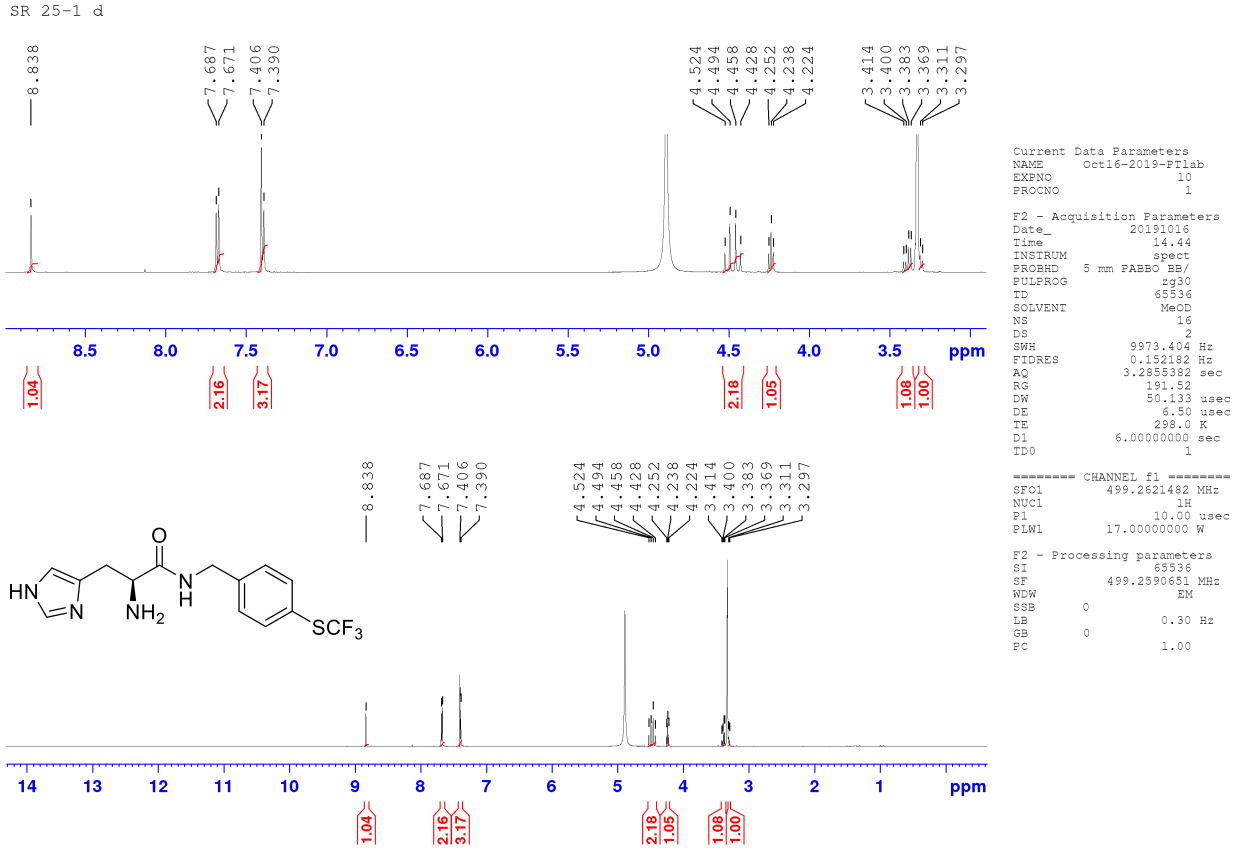
^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethoxy)benzyl)propanamide **(5q):**



^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethoxy)benzyl)propanamide **(5q):**



^{1}H NMR (500 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-((trifluoromethyl)thio)benzyl)propanamide **(5r):**



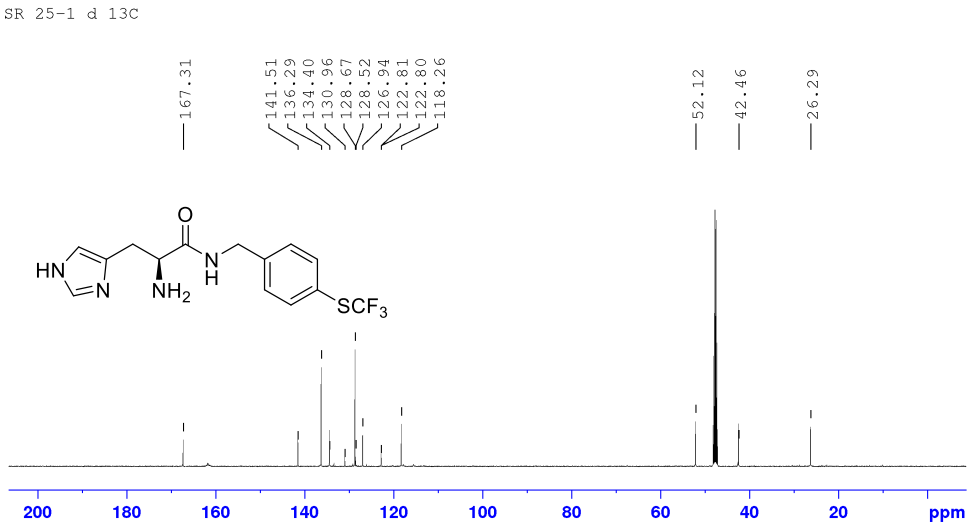

```
Current Data Parameters
NAME      Oct16-2019-PTlab
EXPNO                   10
PROCNO                   1

F2 - Acquisition Parameters
Date_              20191016
Time                 14.44
INSTRUM              spect
PROBHD    5 mm PABBO BB/
PULPROG              zg30
TD                  65536
SOLVENT              MeOD
NS                     16
DS                      2
SWH              9973.404 Hz
FIDRES           0.152182 Hz
AQ               3.2855382 sec
RG                  191.52
DW                 50.133 usec
DE                  6.50 usec
TE                 298.0 K
D1              6.00000000 sec
TD0                     1

======== CHANNEL f1 ========
SFO1         499.2621482 MHz
NUC1                   1H
P1                  10.00 usec
PLW1        17.00000000 W

F2 - Processing parameters
SI                  65536
SF           499.2590651 MHz
WDW                    EM
SSB       0
LB                  0.30 Hz
GB        0
PC                  1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-((trifluoromethyl)thio)benzyl)propanamide **(5r):**




```
Current Data Parameters
NAME      Nov17-2019-PTlab
EXPNO                   50
PROCNO                   1

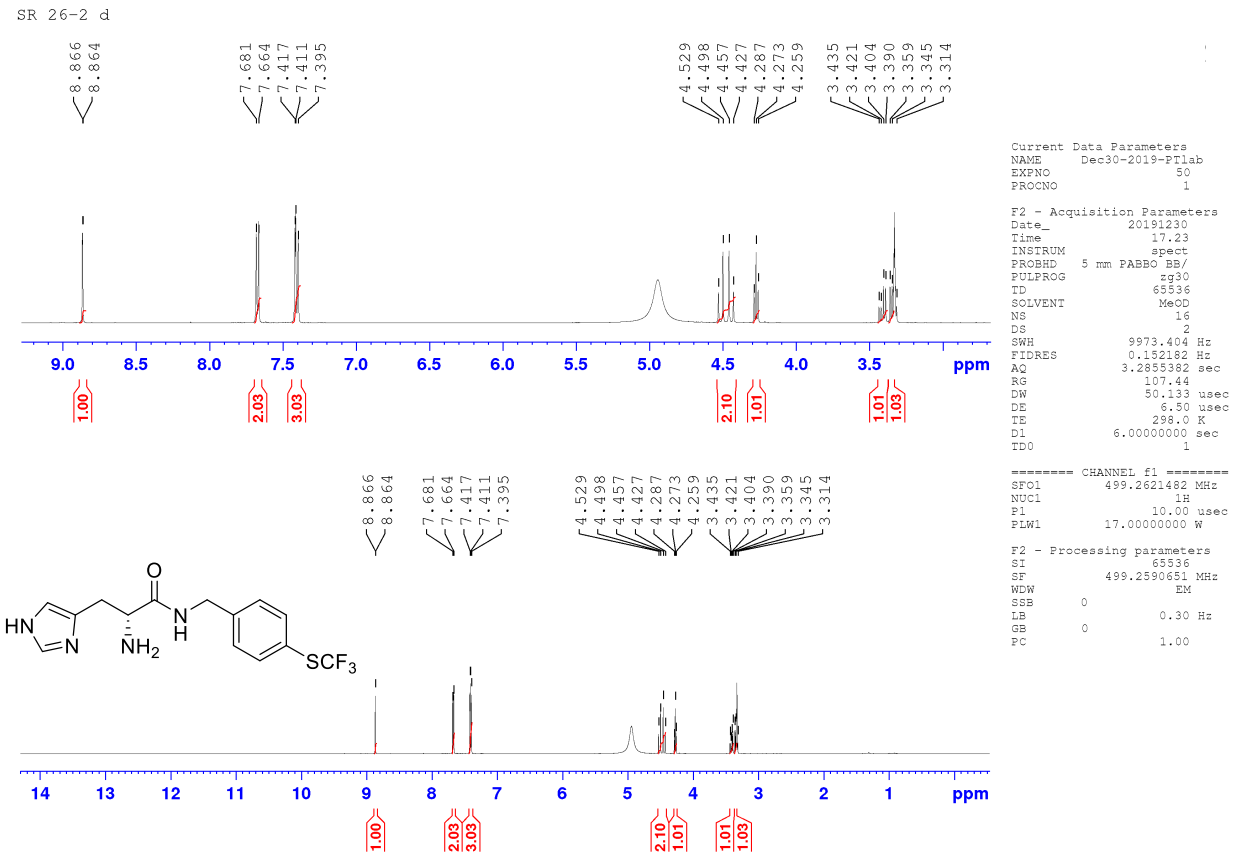
F2 - Acquisition Parameters
Date_           20191117
Time              19.18
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG         zgpg30
TD               65536
SOLVENT           MeOD
NS                 512
DS                   4
SWH           29761.904 Hz
FIDRES         0.454131 Hz
AQ            1.1010048 sec
RG               191.52
DW               16.800 usec
DE               10.00 usec
TE               300.0 K
D1            2.00000000 sec
D11           0.03000000 sec
TD0                  1

======== CHANNEL f1 ========
SFO1         125.5513451 MHz
NUC1               13C
P1               10.00 usec
PLW1        71.00000000 W

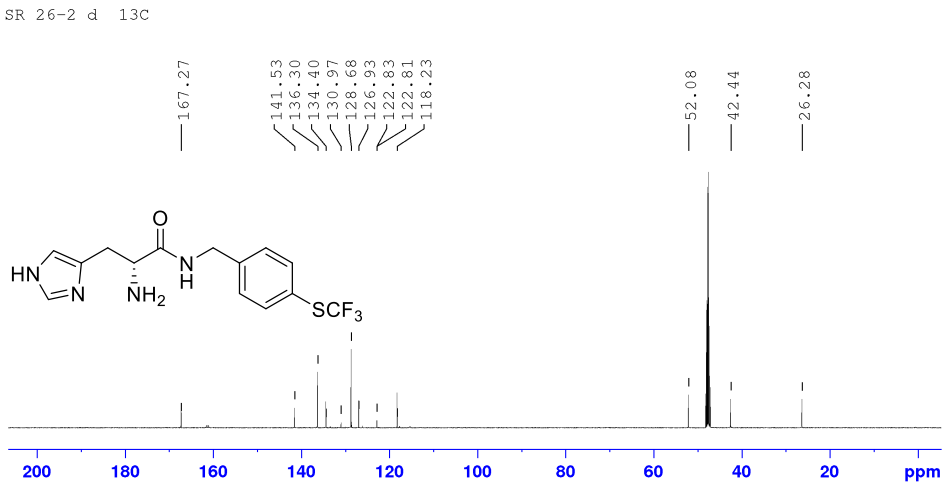
======== CHANNEL f2 ========
SFO2         499.2610621 MHz
NUC2               1H
CPDPRG[2         waltz16
PCPD2             80.00 usec
PLW2        17.00000000 W
PLW12        0.26563001 W
PLW13        0.17000000 W

F2 - Processing parameters
SI                32768
SF           125.5387918 MHz
WDW                EM
SSB       0
LB                1.00 Hz
GB        0
PC                1.40
```

^{1}H NMR (500 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-((trifluoromethyl)thio)benzyl)propanamide **(5s):**



^{13}C NMR (126 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-((trifluoromethyl)thio)benzyl)propanamide **(5s):**




```
Current Data Parameters
NAME      Dec30-2019-PTlab
EXPNO                   60
PROCNO                   1

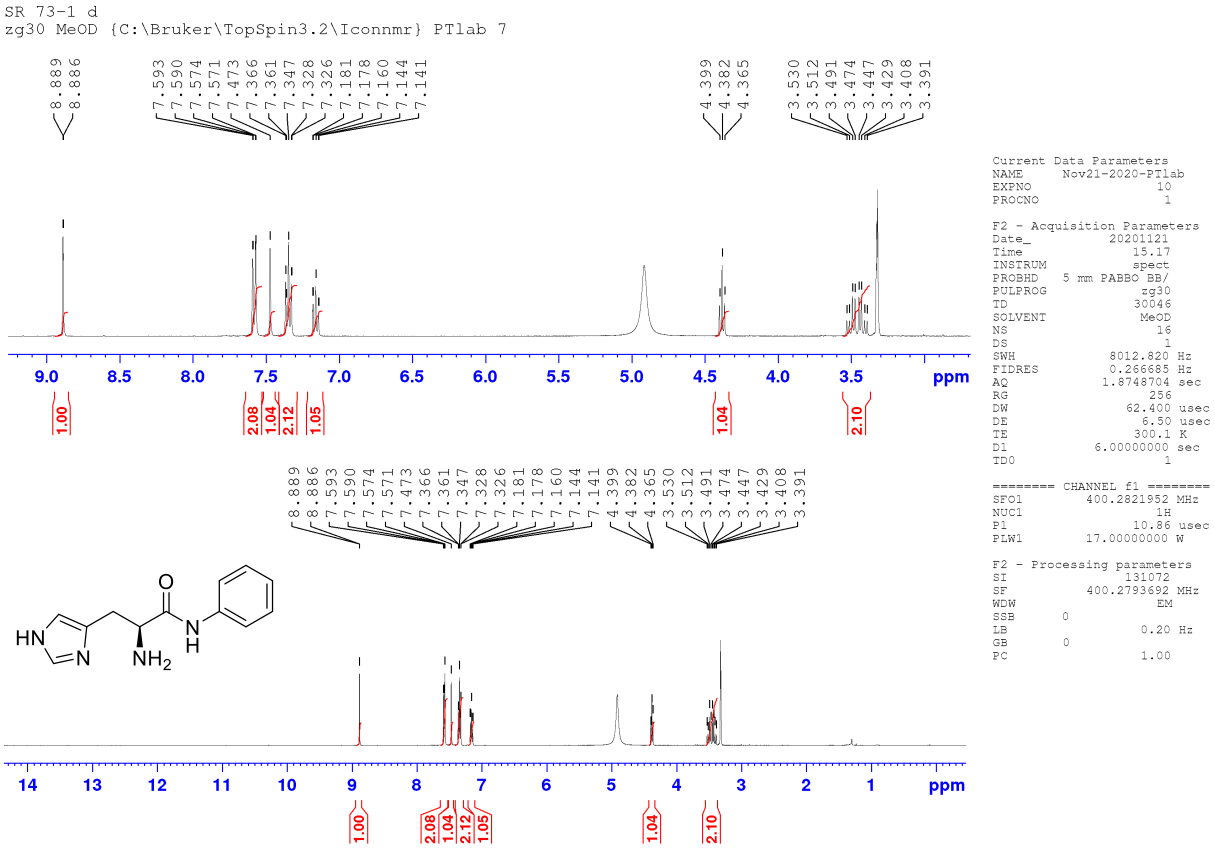
F2 - Acquisition Parameters
Date_            20191230
Time                18.22
INSTRUM              spect
PROBHD   5 mm PABBO BB/
PULPROG           zgpg30
TD                 65536
SOLVENT             MeOD
NS                   512
DS                     4
SWH            29761.904 Hz
FIDRES         0.454131 Hz
AQ             1.1010048 sec
RG                191.52
DW                16.800 usec
DE                10.00 usec
TE                300.0 K
D1             2.00000000 sec
D11            0.03000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1         125.5513451 MHz
NUC1                  13C
P1                 10.00 usec
PLW1         71.00000000 W

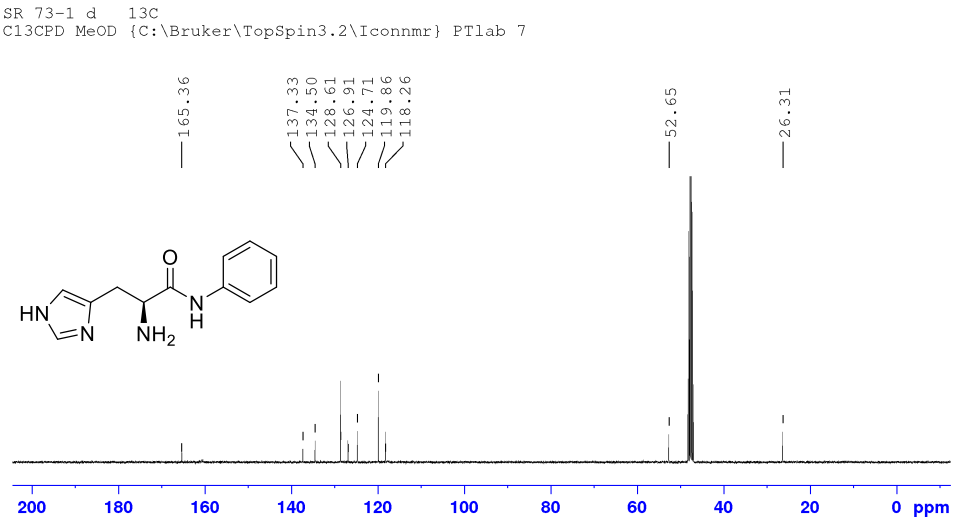
======== CHANNEL f2 ========
SFO2         499.2610621 MHz
NUC2                  1H
CPDPRG[2          waltz16
PCPD2              80.00 usec
PLW2         17.00000000 W
PLW12         0.26563001 W
PLW13         0.17000000 W

F2 - Processing parameters
SI                  32768
SF           125.5387918 MHz
WDW                    EM
SSB       0
LB                 1.00 Hz
GB        0
PC                 1.40
```

^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenylpropanamide **(6a):**



^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenylpropanamide **(6a):**




```
Current Data Parameters
NAME      Nov21-2020-PTlab
EXPNO                   20
PROCNO                   1

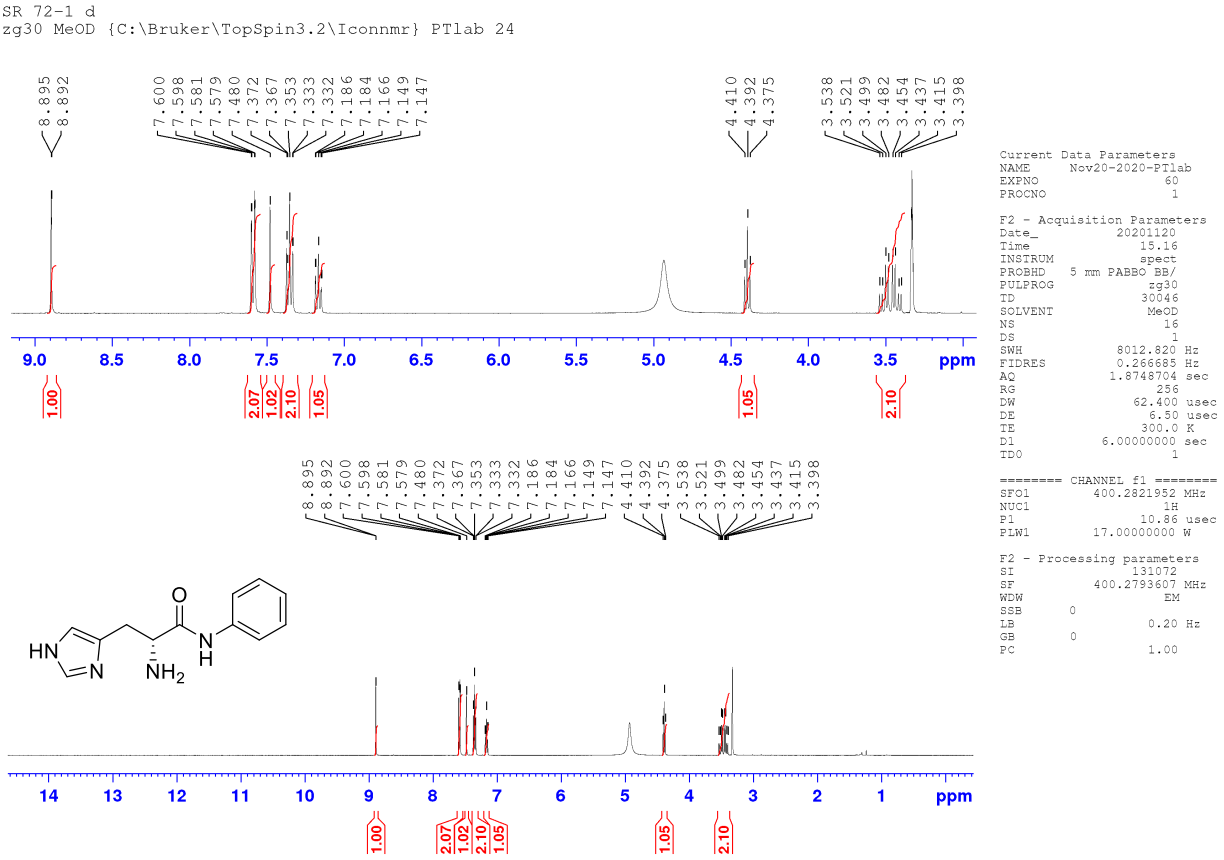
F2 - Acquisition Parameters
Date_           20201121
Time              15.58
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG          zgpg30
TD                65536
SOLVENT             MeOD
NS                  512
DS                    4
SWH           24038.461 Hz
FIDRES         0.366798 Hz
AQ             1.3631488 sec
RG                 2050
DW               20.800 usec
DE                6.50 usec
TE                300.0 K
D1            2.00000000 sec
D11           0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1         100.6605506 MHz
NUC1               13C
P1               11.00 usec
PLW1         48.00000000 W

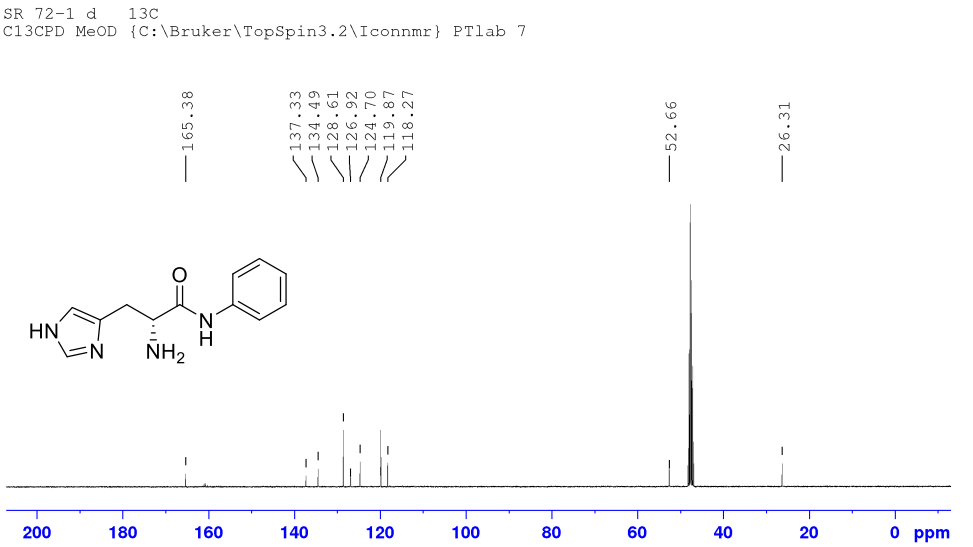
======== CHANNEL f2 ========
SFO2         400.2816011 MHz
NUC2               1H
CPDPRG[2         waltz16
PCPD2             90.00 usec
PLW2         17.00000000 W
PLW12         0.24753000 W
PLW13         0.20050000 W

F2 - Processing parameters
SI                 32768
SF           100.6504861 MHz
WDW                  EM
SSB       0
LB                1.00 Hz
GB        0
PC                1.40
```

^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenylpropanamide **(6b):**



^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenylpropanamide **(6b):**




```
Current Data Parameters
NAME      Nov20-2020-PTlab
EXPNO                  100
PROCNO                   1

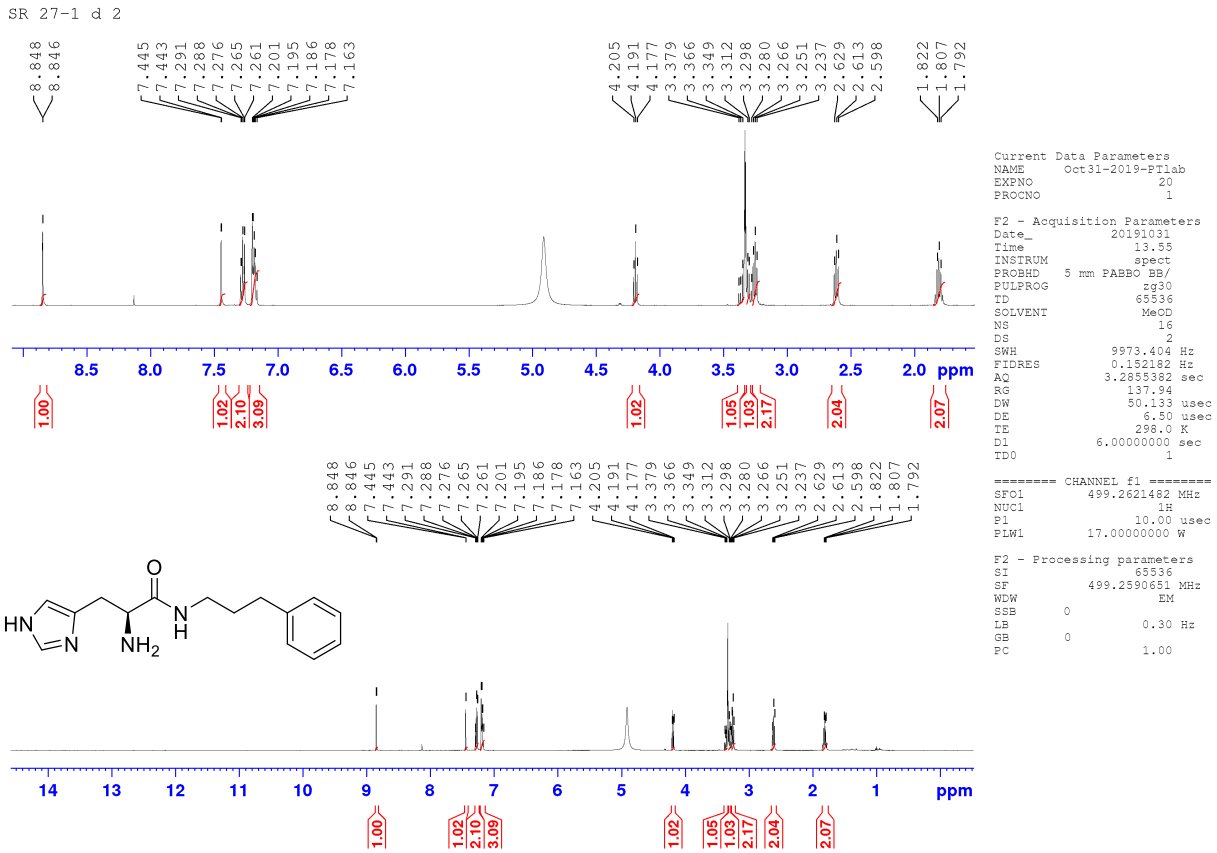
F2 - Acquisition Parameters
Date_            20201120
Time                17.54
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG           zgpg30
TD                 65536
SOLVENT             MeOD
NS                   512
DS                     4
SWH            24038.461 Hz
FIDRES          0.366798 Hz
AQ             1.3631488 sec
RG                  2050
DW                20.800 usec
DE                 6.50 usec
TE                300.0 K
D1           2.00000000 sec
D11          0.03000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1         100.6605506 MHz
NUC1                  13C
P1                 11.00 usec
PLW1         48.00000000 W

======== CHANNEL f2 ========
SFO2         400.2816011 MHz
NUC2                  1H
CPDPRG[2          waltz16
PCPD2               90.00 usec
PLW2         17.00000000 W
PLW12         0.24753000 W
PLW13         0.20050000 W

F2 - Processing parameters
SI                   32768
SF            100.6504861 MHz
WDW                    EM
SSB       0
LB                  1.00 Hz
GB        0
PC                  1.40
```

^{1}H NMR (500 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-phenylpropyl)propanamide **(7a):**

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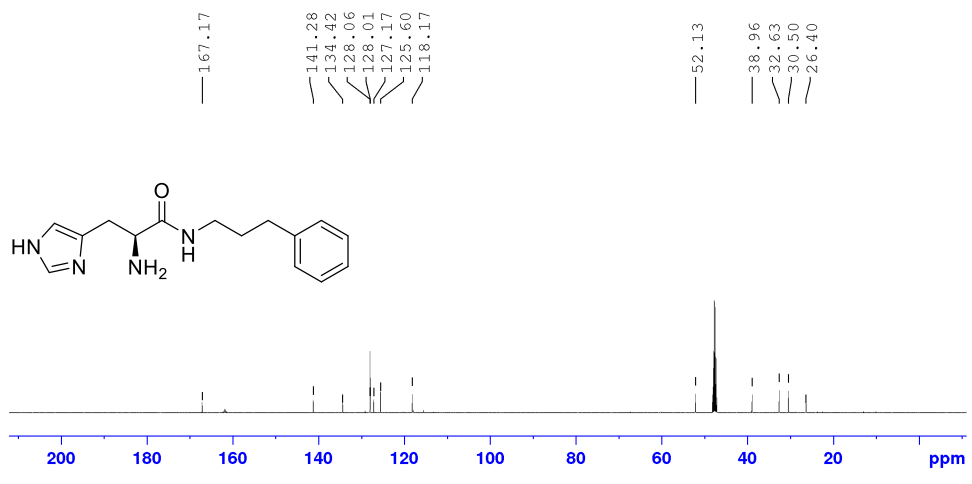
Current Data Parameters
NAME      Oct31-2019-PTlab
EXPNO                   20
PROCNO                   1

F2 - Acquisition Parameters
Date_            20191031
Time                13.55
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG             zg30
TD                 65536
SOLVENT             MeOD
NS                    16
DS                     2
SWH              9973.404 Hz
FIDRES           0.152182 Hz
AQ               3.2855382 sec
RG                 137.94
DW                  50.133 usec
DE                   6.50 usec
TE                 298.0 K
D1              6.00000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1          499.2621482 MHz
NUC1                  1H
P1                 10.00 usec
PLW1          17.00000000 W

F2 - Processing parameters
SI                 65536
SF            499.2590651 MHz
WDW                   EM
SSB       0
LB                 0.30 Hz
GB        0
PC                 1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-phenylpropyl)propanamide **(7a):**




```
Current Data Parameters
NAME       Nov17-2019-PTlab
EXPNO                    60
PROCNO                    1

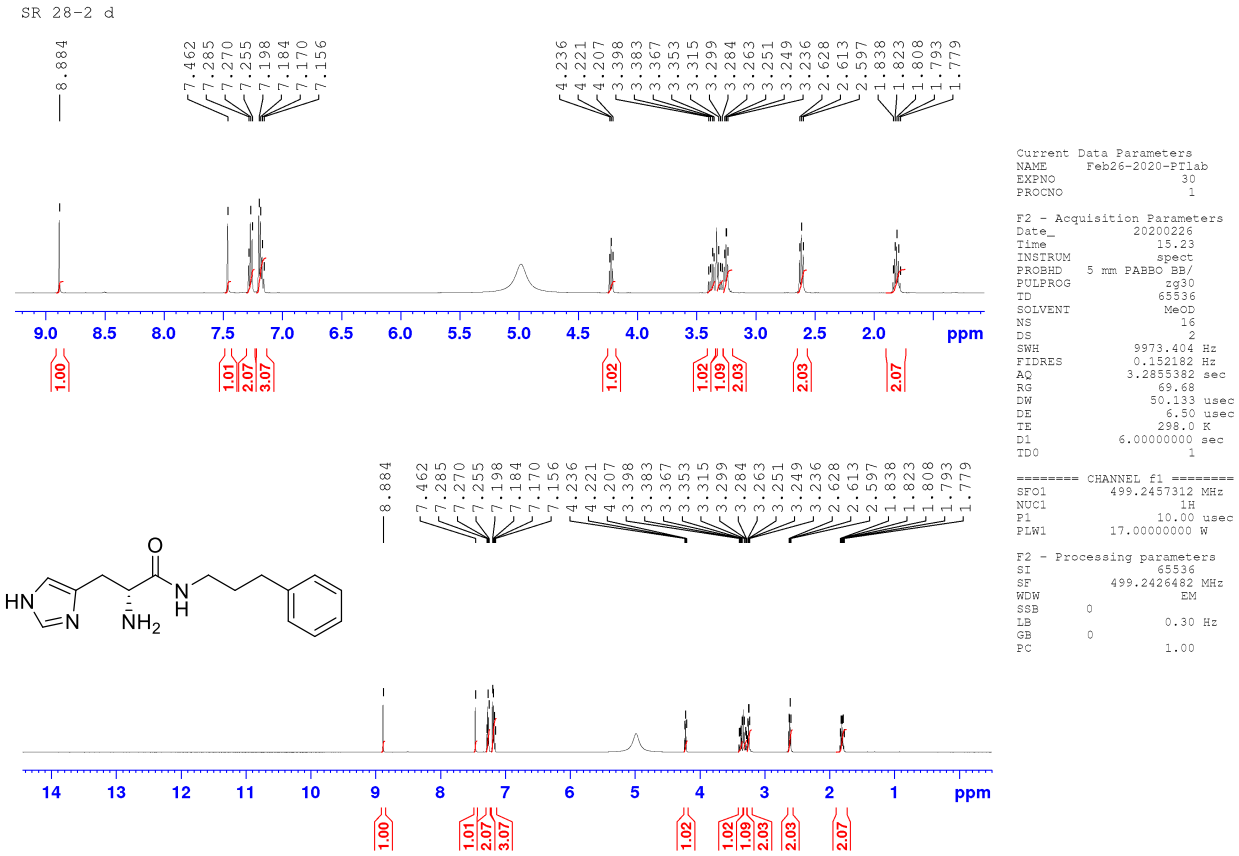
F2 - Acquisition Parameters
Date_               20191117
Time                   19.49
INSTRUM                spect
PROBHD    5 mm PABBO BB/
PULPROG             zgpg30
TD                   65536
SOLVENT                MeOD
NS                       512
DS                         4
SWH                29761.904 Hz
FIDRES             0.454131 Hz
AQ                 1.1010048 sec
RG                    191.52
DW                  16.800 usec
DE                  10.00 usec
TE                    300.0 K
D1              2.00000000 sec
D11             0.03000000 sec
TD0                       1

======== CHANNEL f1 ========
SFO1           125.5513451 MHz
NUC1                     13C
P1                    10.00 usec
PLW1           71.00000000 W

======== CHANNEL f2 ========
SFO2           499.2610621 MHz
NUC2                     1H
CPDPRG[2          waltz16
PCPD2                80.00 usec
PLW2           17.00000000 W
PLW12           0.26563001 W
PLW13           0.17000000 W

F2 - Processing parameters
SI                    32768
SF           125.5387918 MHz
WDW                      EM
SSB       0
LB                    1.00 Hz
GB        0
PC                    1.40
```

^{1}H NMR (500 MHz, MeOH-d$_4$), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-phenylpropyl)propanamide **(7b):**



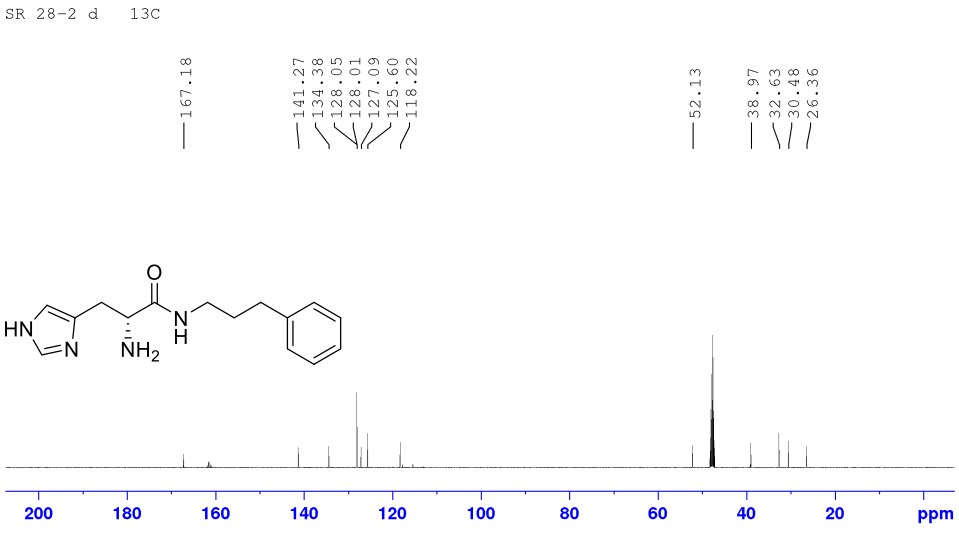

```
Current Data Parameters
NAME       Feb26-2020-PTlab
EXPNO                   30
PROCNO                   1

F2 - Acquisition Parameters
Date_               20200226
Time                   15.23
INSTRUM                spect
PROBHD    5 mm PABBO BB/
PULPROG                zg30
TD                    65536
SOLVENT                MeOD
NS                       16
DS                        2
SWH                9973.404 Hz
FIDRES            0.152182 Hz
AQ                3.2855382 sec
RG                     69.68
DW                  50.133 usec
DE                   6.50 usec
TE                   298.0 K
D1                6.00000000 sec
TD0                      1

======== CHANNEL f1 ========
SFO1            499.2457312 MHz
NUC1                    1H
P1                   10.00 usec
PLW1           17.00000000 W

F2 - Processing parameters
SI                    65536
SF              499.2426482 MHz
WDW                     EM
SSB       0
LB                    0.30 Hz
GB        0
PC                    1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-phenylpropyl)propanamide **(7b):**




```
Current Data Parameters
NAME      Feb26-2020-PTlab
EXPNO                   40
PROCNO                   1

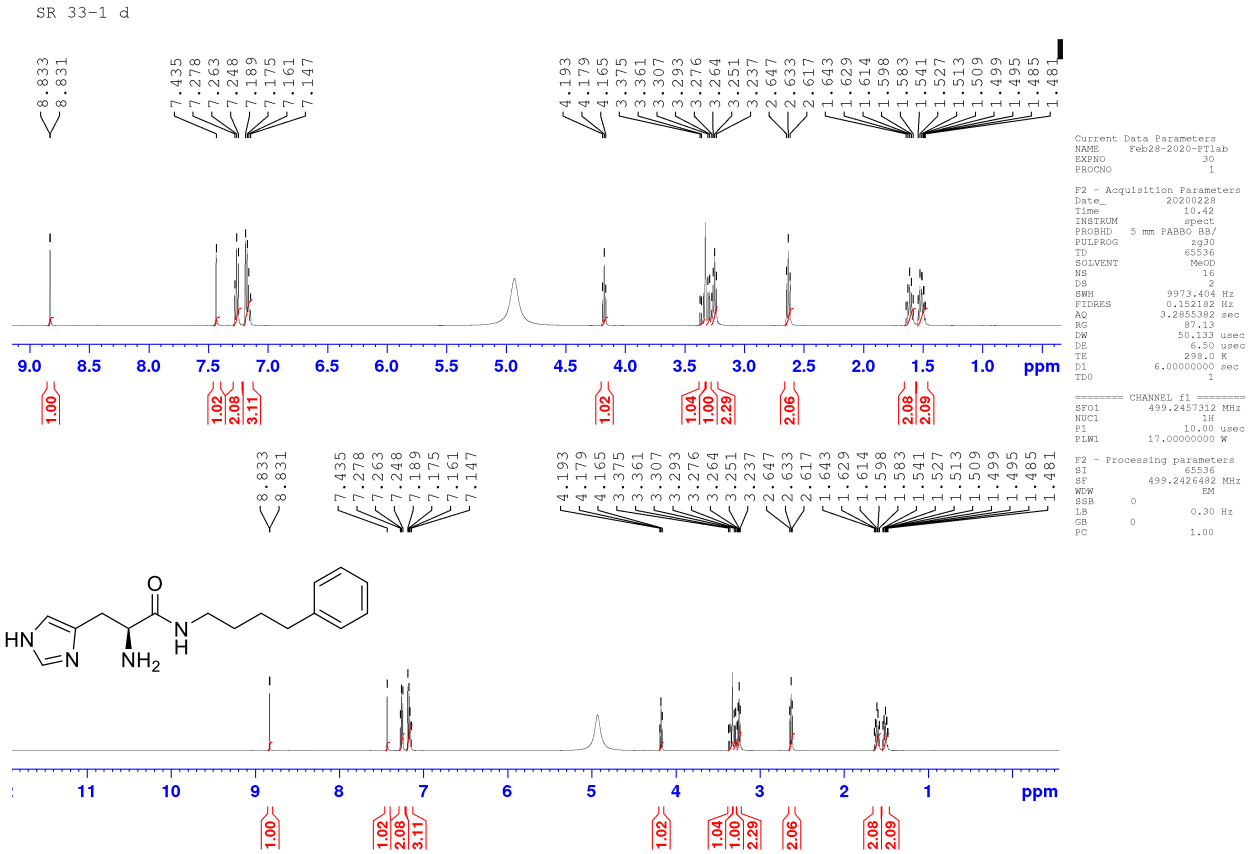
F2 - Acquisition Parameters
Date_              20200226
Time                 16.34
INSTRUM              spect
PROBHD   5 mm PABBO BB/
PULPROG           zgpg30
TD                 65536
SOLVENT              MeOD
NS                     512
DS                       4
SWH            29761.904 Hz
FIDRES          0.454131 Hz
AQ             1.1010048 sec
RG                  191.52
DW                16.800 usec
DE                10.00 usec
TE                300.0 K
D1             2.00000000 sec
D11            0.03000000 sec
TD0                      1

======== CHANNEL f1 ========
SFO1         125.5472166 MHz
NUC1                   13C
P1                   10.00 usec
PLW1         71.00000000 W

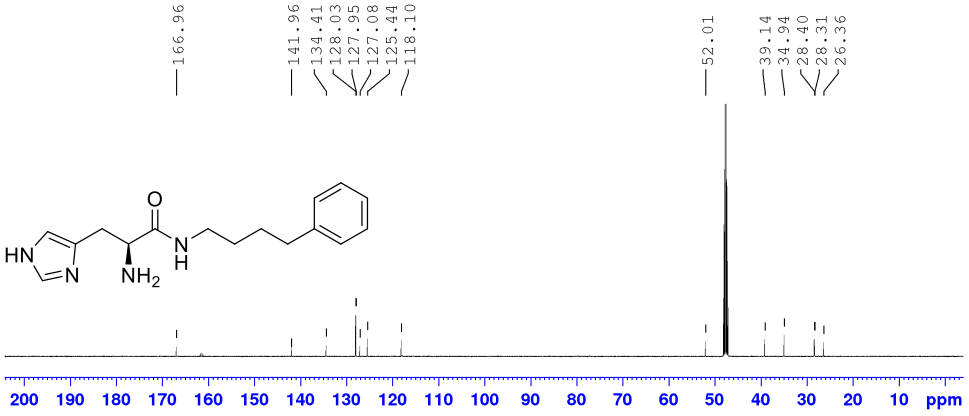
======== CHANNEL f2 ========
SFO2         499.2446452 MHz
NUC2                   1H
CPDPRG[2          waltz16
PCPD2                80.00 usec
PLW2         17.00000000 W
PLW12         0.26563001 W
PLW13         0.17000000 W

F2 - Processing parameters
SI                   32768
SF           125.5346637 MHz
WDW                     EM
SSB       0
LB                    1.00 Hz
GB        0
PC                    1.40
```

^{1}H NMR (500 MHz, MeOH-d$_4$), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-phenylbutyl)propanamide **(8a):**



^{13}C NMR (126 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-phenylbutyl)propanamide **(8a):**




```
Current Data Parameters
NAME      Feb28-2020-PTlab
EXPNO                   40
PROCNO                   1

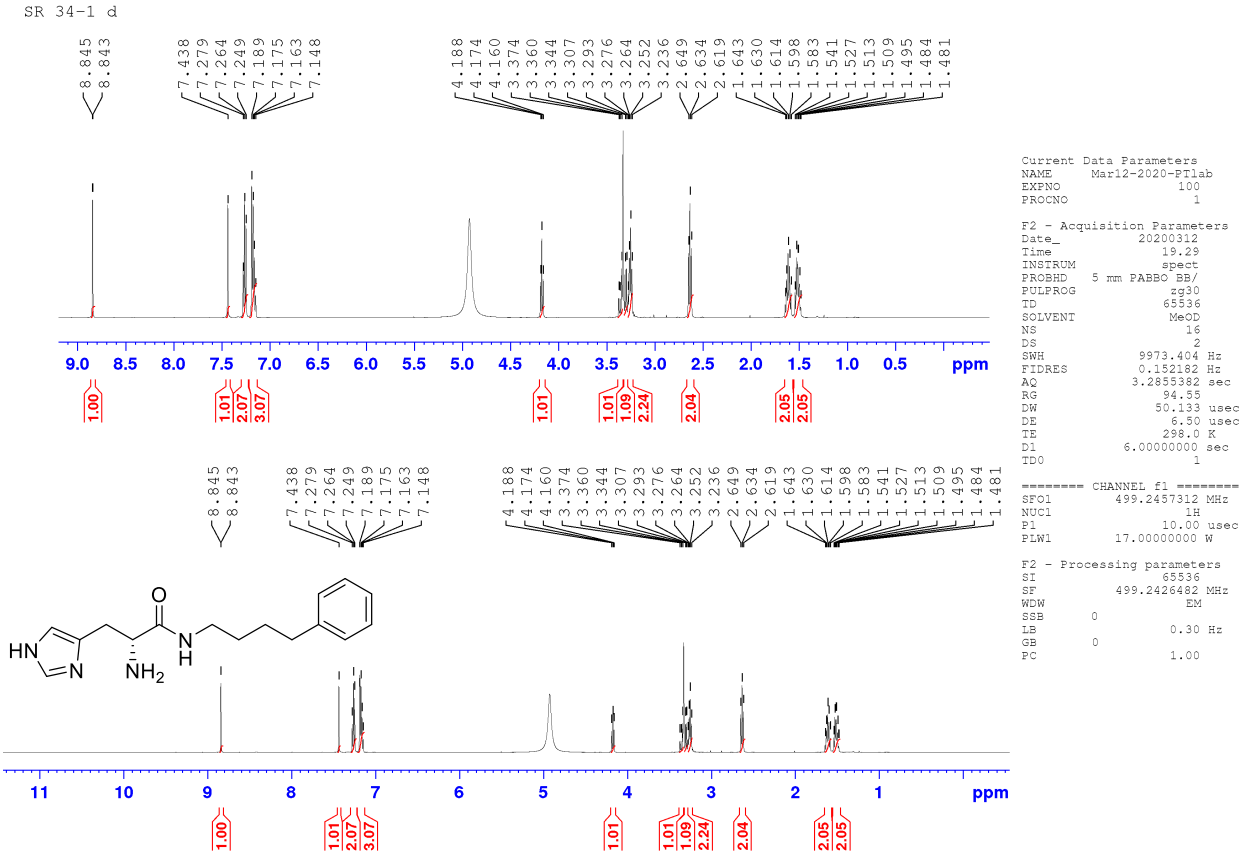
F2 - Acquisition Parameters
Date_              20200228
Time                 11.46
INSTRUM              spect
PROBHD    5 mm PABBO BB/
PULPROG            zgpg30
TD                  65536
SOLVENT               MeOD
NS                     512
DS                       4
SWH              29761.904 Hz
FIDRES           0.454131 Hz
AQ               1.1010048 sec
RG                  191.52
DW                 16.800 usec
DE                 10.00 usec
TE                 300.0 K
D1              2.00000000 sec
D11             0.03000000 sec
TD0                      1

======== CHANNEL f1 ========
SFO1          125.5472166 MHz
NUC1                    13C
P1                   10.00 usec
PLW1          71.00000000 W

======== CHANNEL f2 ========
SFO2          499.2446452 MHz
NUC2                    1H
CPDPRG[2           waltz16
PCPD2                80.00 usec
PLW2          17.00000000 W
PLW12          0.26563001 W
PLW13          0.17000000 W

F2 - Processing parameters
SI                    32768
SF            125.5346637 MHz
WDW                      EM
SSB       0
LB                    1.00 Hz
GB        0
PC                    1.40
```

^{1}H NMR (500 MHz, MeOH-d$_4$), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-phenylbutyl)propanamide **(8b):**




```
Current Data Parameters
NAME      Mar12-2020-PTlab
EXPNO                  100
PROCNO                   1

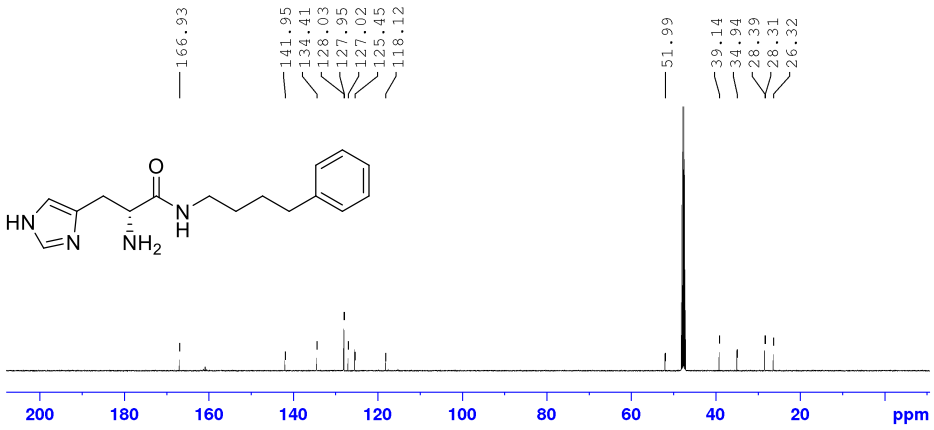
F2 - Acquisition Parameters
Date_            20200312
Time                19.29
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG             zg30
TD                 65536
SOLVENT             MeOD
NS                    16
DS                     2
SWH              9973.404 Hz
FIDRES           0.152182 Hz
AQ               3.2855382 sec
RG                 94.55
DW                 50.133 usec
DE                  6.50 usec
TE                298.0 K
D1             6.00000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1          499.2457312 MHz
NUC1                   1H
P1                  10.00 usec
PLW1          17.00000000 W

F2 - Processing parameters
SI                 65536
SF            499.2426482 MHz
WDW                   EM
SSB       0
LB                  0.30 Hz
GB        0
PC                  1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-phenylbutyl)propanamide **(8b):**

SR 34-1 d 13C




```
Current Data Parameters
NAME      Mar12-2020-PTlab
EXPNO                  110
PROCNO                   1

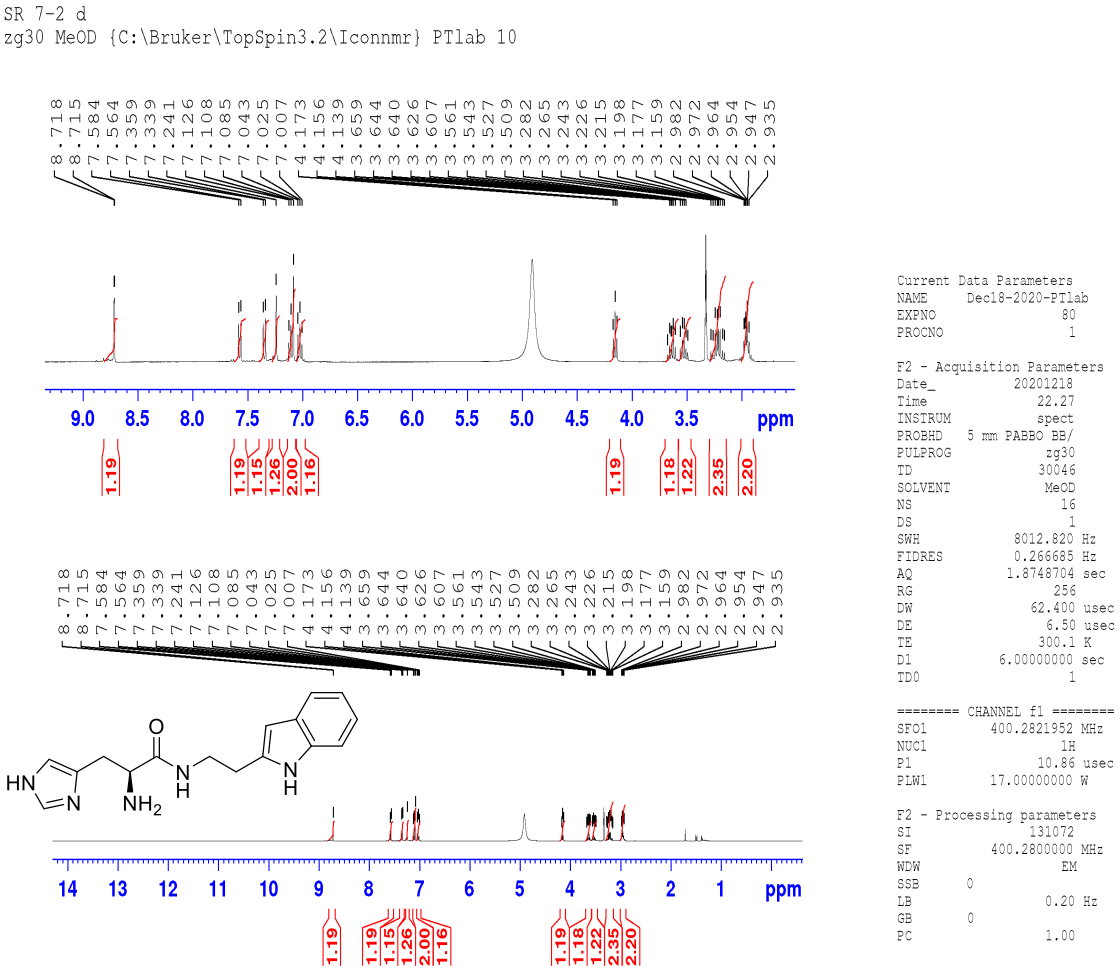
F2 - Acquisition Parameters
Date_           20200312
Time              20.05
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG         zgpg30
TD                65536
SOLVENT             MeOD
NS                    512
DS                      4
SWH           29761.904 Hz
FIDRES         0.454131 Hz
AQ            1.1010048 sec
RG               191.52
DW               16.800 usec
DE               10.00 usec
TE                300.0 K
D1           2.00000000 sec
D11          0.03000000 sec
TD0                   1

======== CHANNEL f1 ========
SFO1        125.5472166 MHz
NUC1               13C
P1               10.00 usec
PLW1        71.00000000 W

======== CHANNEL f2 ========
SFO2        499.2446452 MHz
NUC2               1H
CPDPRG[2        waltz16
PCPD2            80.00 usec
PLW2        17.00000000 W
PLW12        0.26563001 W
PLW13        0.17000000 W

F2 - Processing parameters
SI                32768
SF          125.5346637 MHz
WDW                  EM
SSB       0
LB                1.00 Hz
GB        0
PC                1.40
```

^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-*N*-(2-(1*H*-indol-3-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide **(9d):**



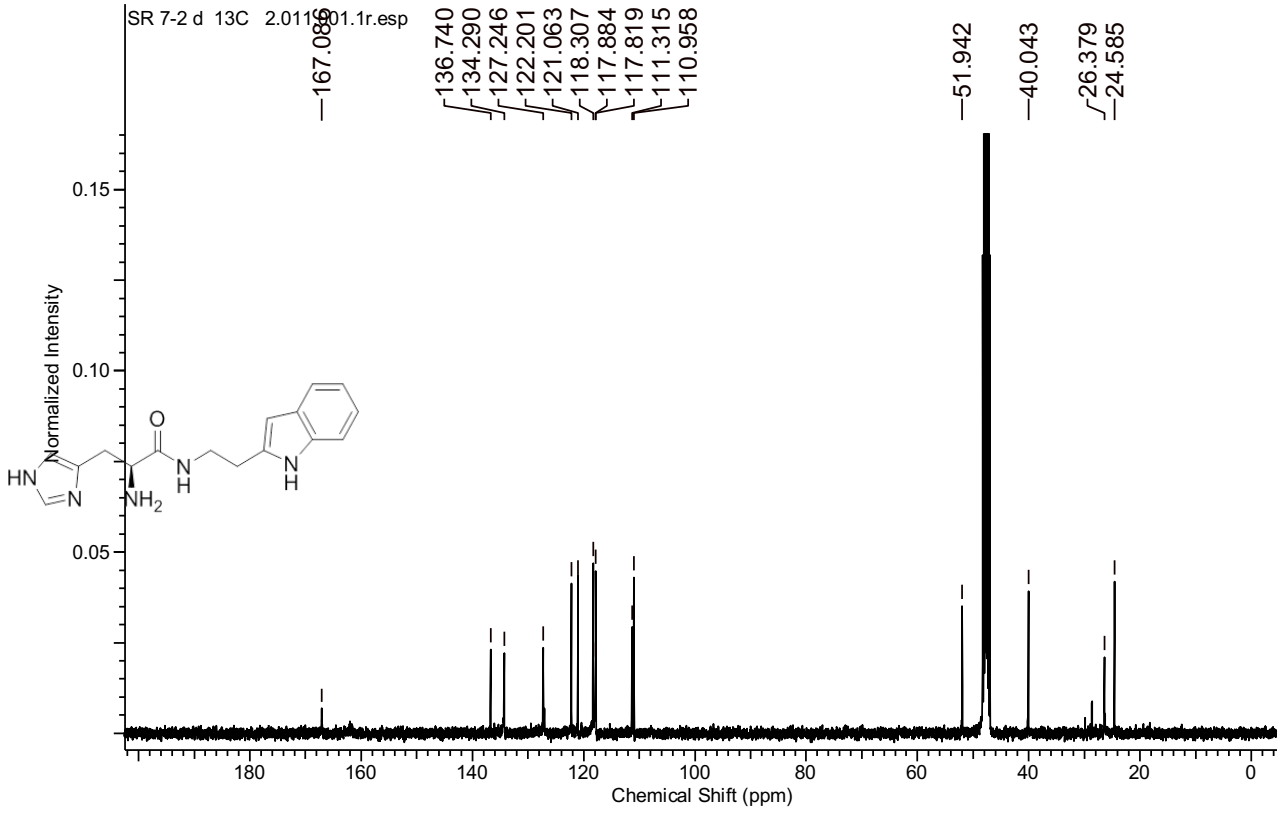

```
Current Data Parameters
NAME      Dec18-2020-PTlab
EXPNO                   80
PROCNO                   1

F2 - Acquisition Parameters
Date_           20201218
Time              22.27
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG             zg30
TD                 30046
SOLVENT             MeOD
NS                    16
DS                     1
SWH            8012.820 Hz
FIDRES         0.266685 Hz
AQ            1.8748704 sec
RG                   256
DW                62.400 usec
DE                 6.50 usec
TE                300.1 K
D1           6.00000000 sec
TD0                    1

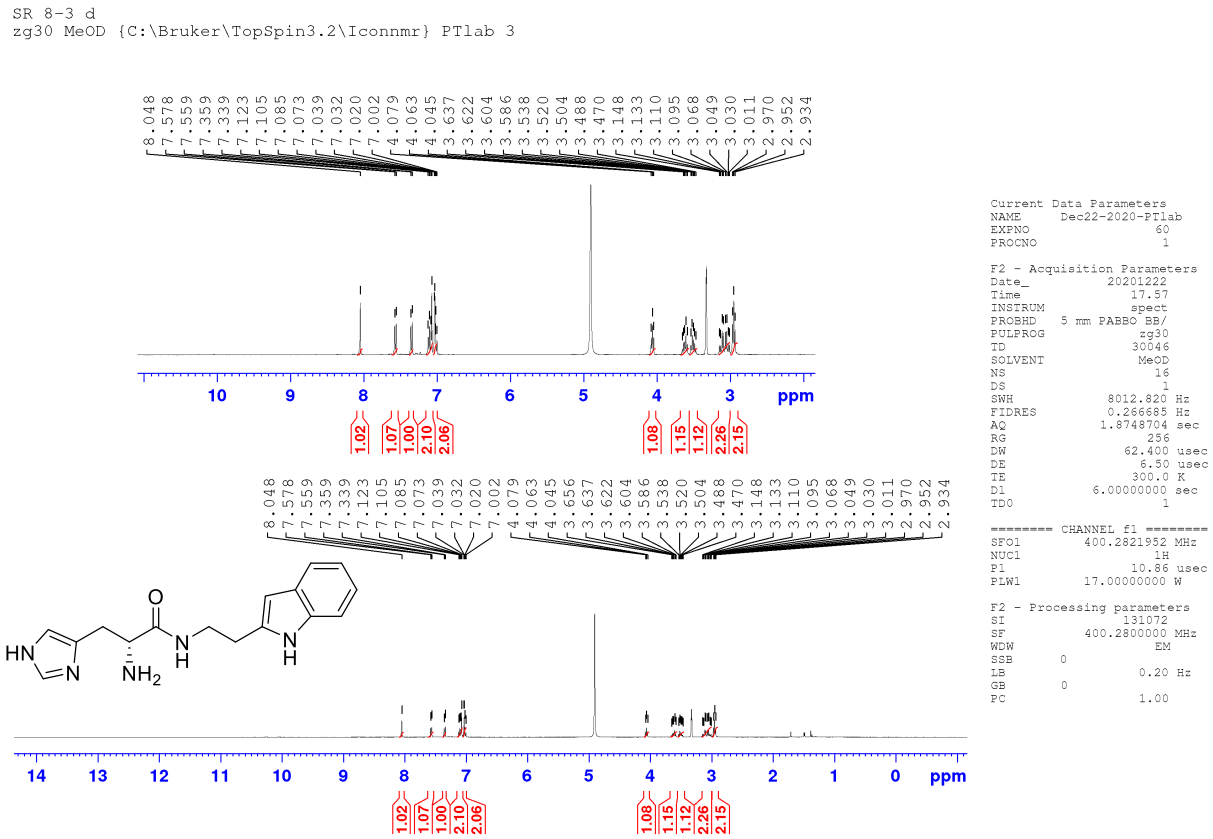
======== CHANNEL f1 ========
SFO1         400.2821952 MHz
NUC1                   1H
P1                 10.86 usec
PLW1        17.00000000 W

F2 - Processing parameters
SI                131072
SF           400.2800000 MHz
WDW                  EM
SSB       0
LB                 0.20 Hz
GB        0
PC                 1.00
```

^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-*N*-(2-(1*H*-indol-3-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide **(9d):**

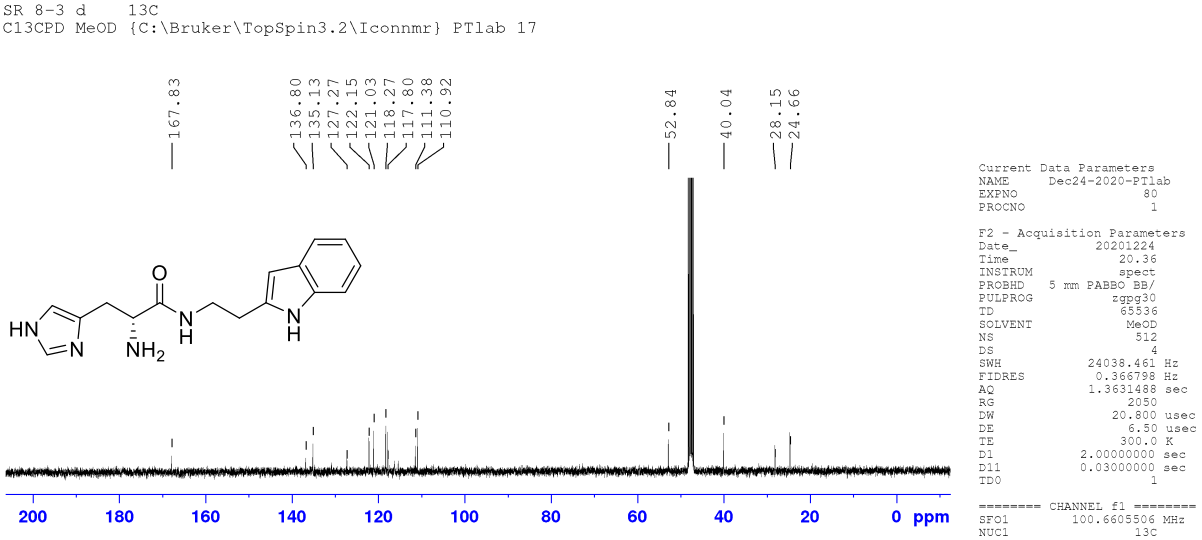


^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-*N*-(2-(1*H*-indol-3-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide **(9e):**



^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-*N*-(2-(1*H*-indol-3-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide

(9e):




```
Current Data Parameters
NAME      Dec24-2020-PTlab
EXPNO                   80
PROCNO                   1

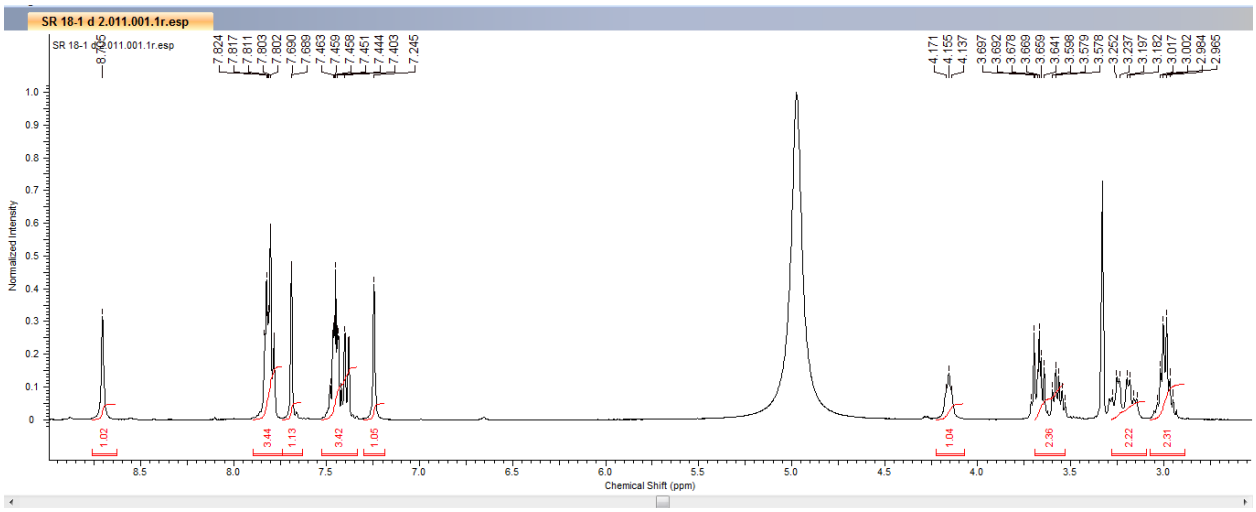
F2 - Acquisition Parameters
Date_           20201224
Time              20.36
INSTRUM           spect
PROBHD   5 mm PABBO BB/
PULPROG         zgpg30
TD               65536
SOLVENT           MeOD
NS                 512
DS                   4
SWH           24038.461 Hz
FIDRES         0.366798 Hz
AQ            1.3631488 sec
RG                2050
DW              20.800 usec
DE               6.50 usec
TE              300.0 K
D1           2.00000000 sec
D11          0.03000000 sec
TD0                  1

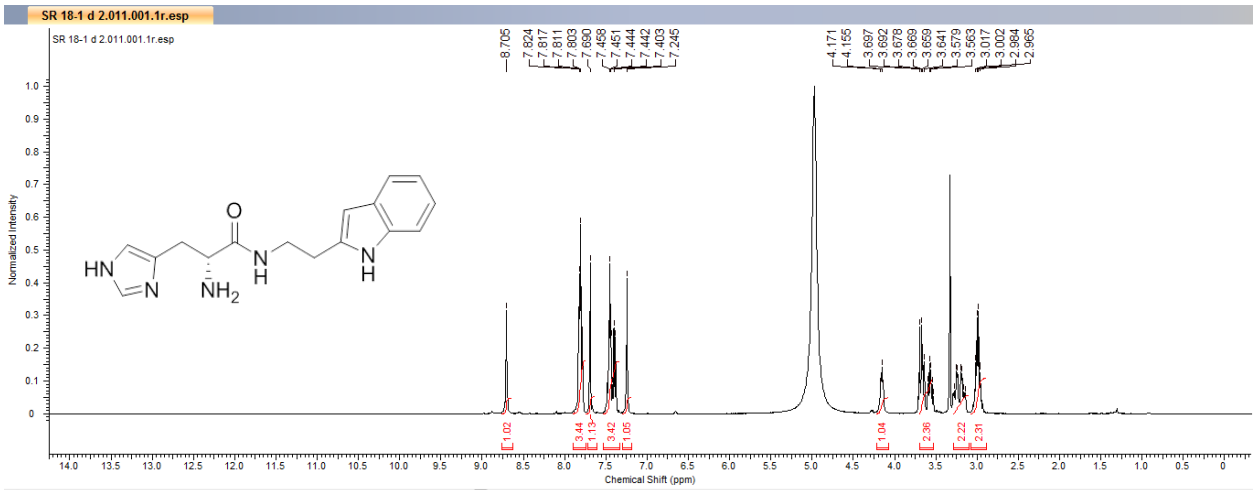
======== CHANNEL f1 ========
SFO1        100.6605506 MHz
NUC1              13C
P1              11.00 usec
PLW1        48.00000000 W

======== CHANNEL f2 ========
SFO2        400.2816011 MHz
NUC2              1H
CPDPRG[2        waltz16
PCPD2           90.00 usec
PLW2        17.00000000 W
PLW12        0.24753000 W
PLW13        0.20050000 W

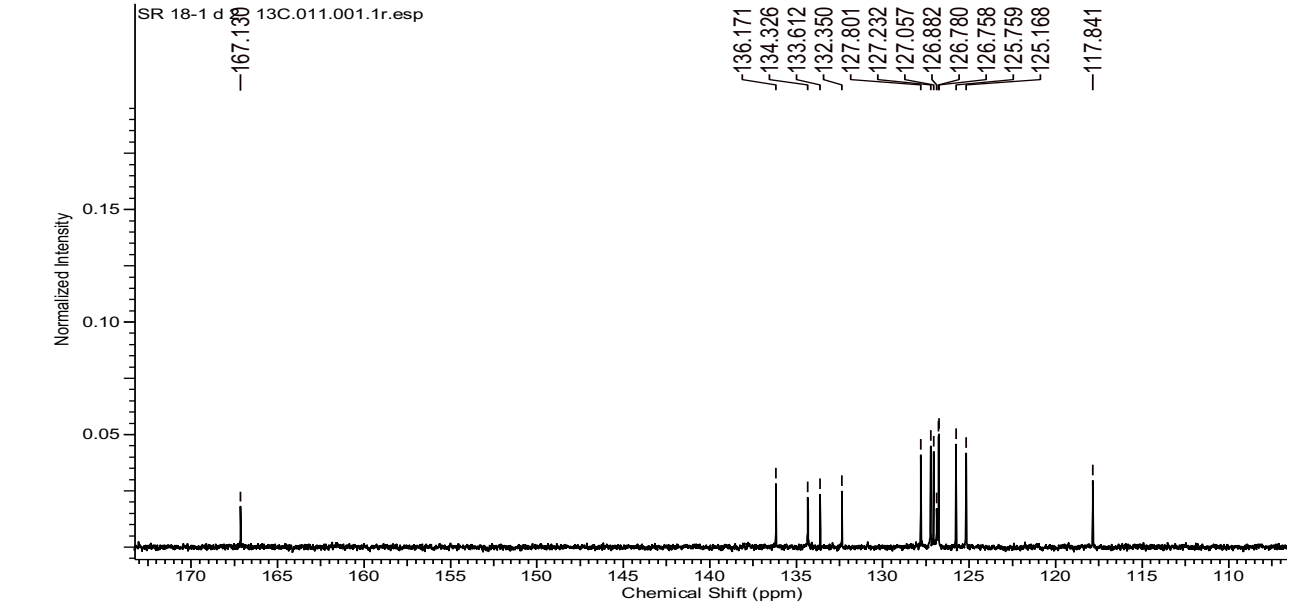
F2 - Processing parameters
SI                32768
SF          100.6504861 MHz
WDW                EM
SSB       0
LB              1.00 Hz
GB        0
PC              1.40
```

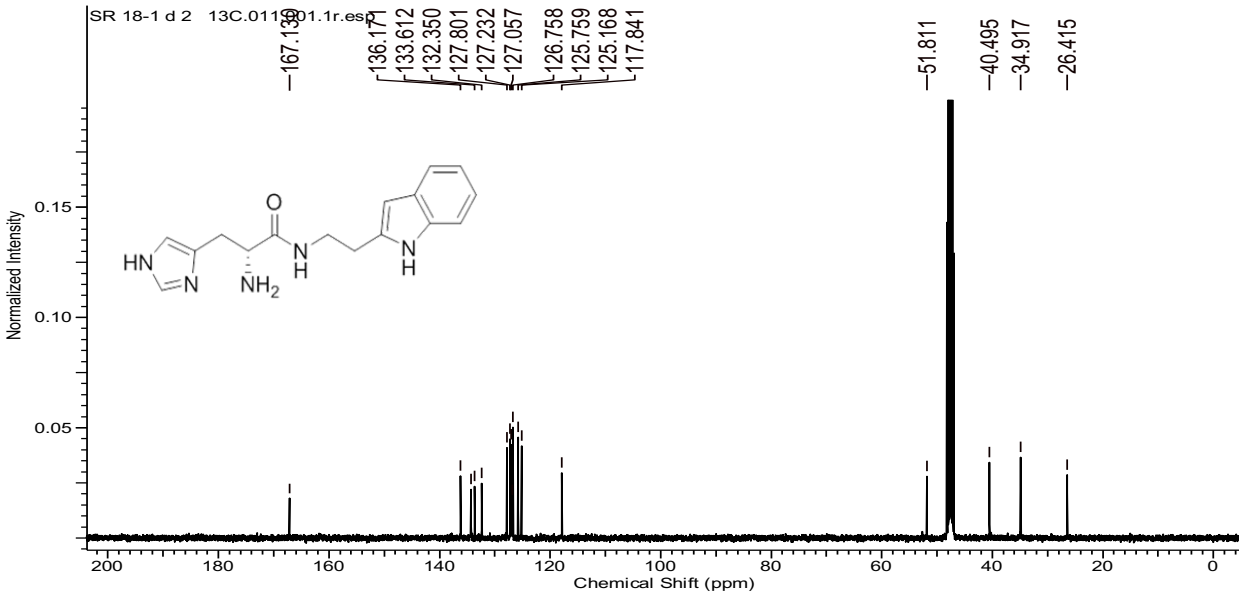
^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(naphthalen-2-yl)ethyl)propanamide **(9f):**



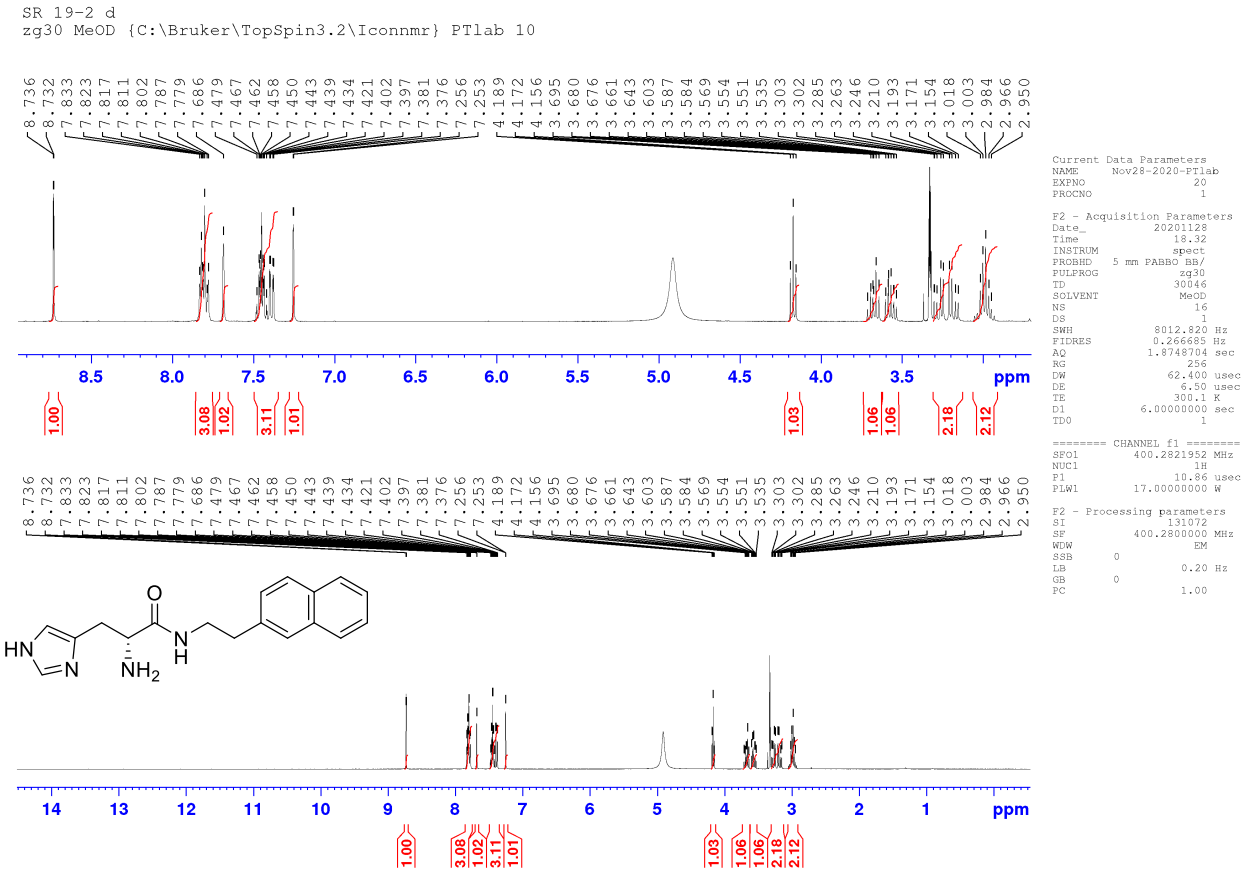


^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(naphthalen-2-yl)ethyl)propanamide **(9f):**

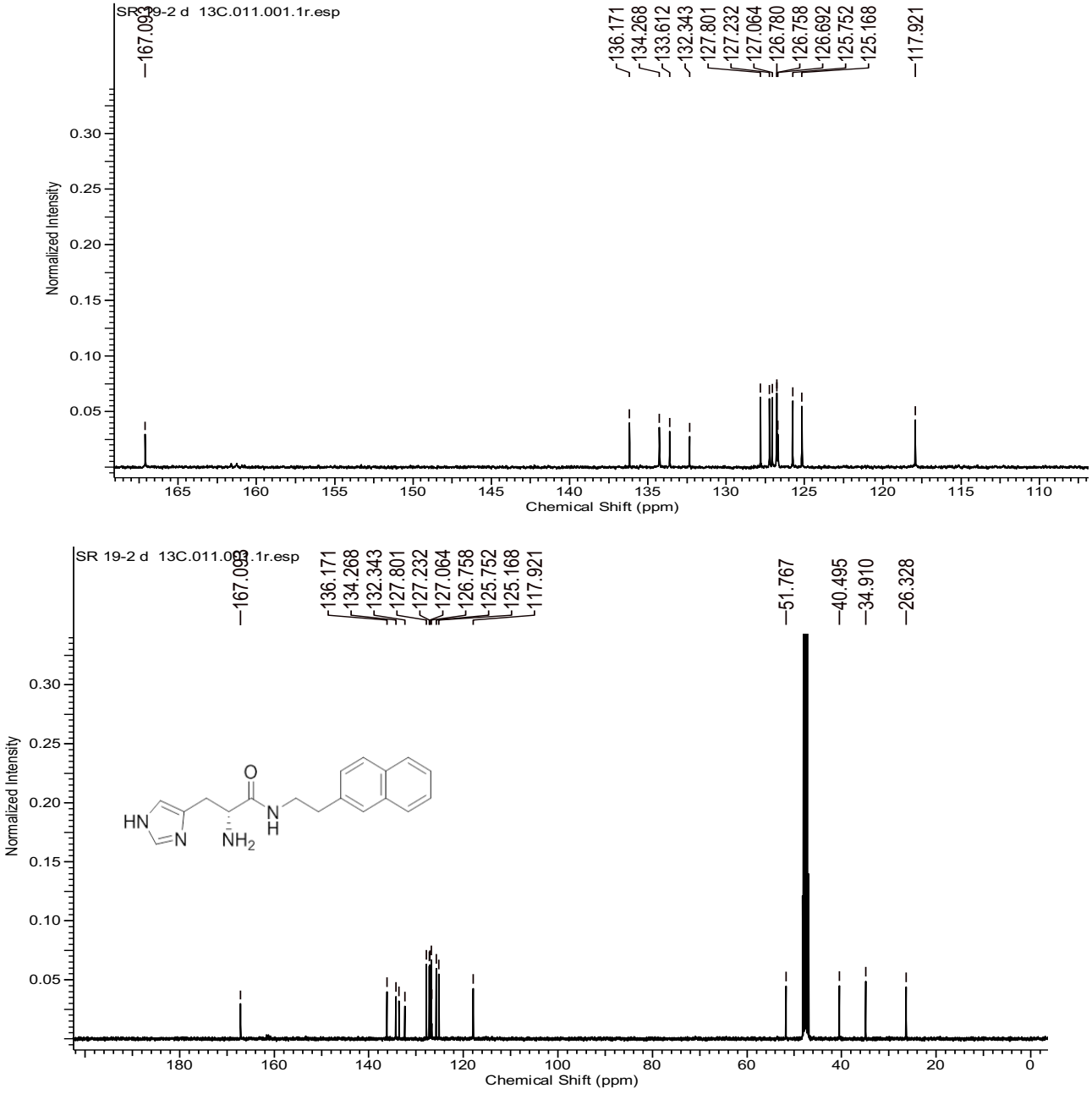




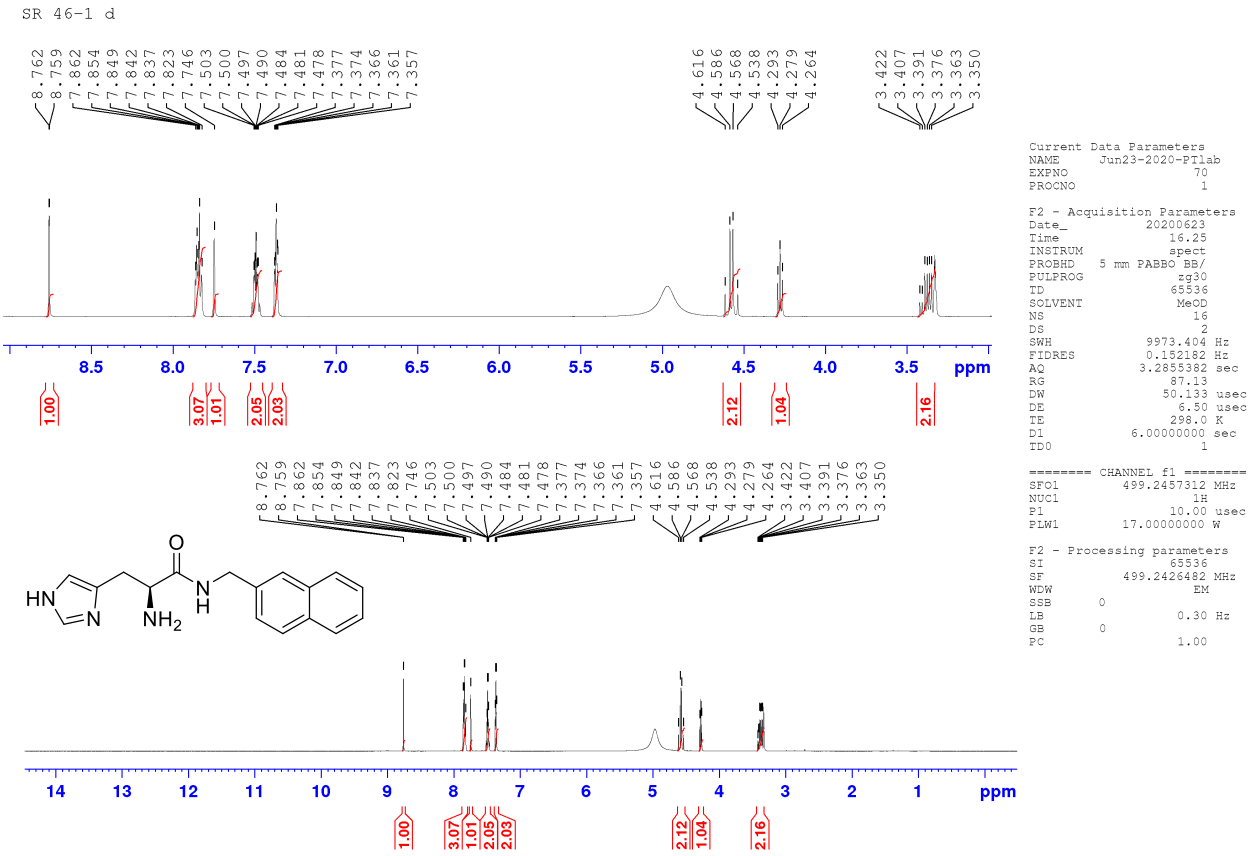
^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(naphthalen-2-yl)ethyl)propanamide **(9g):**



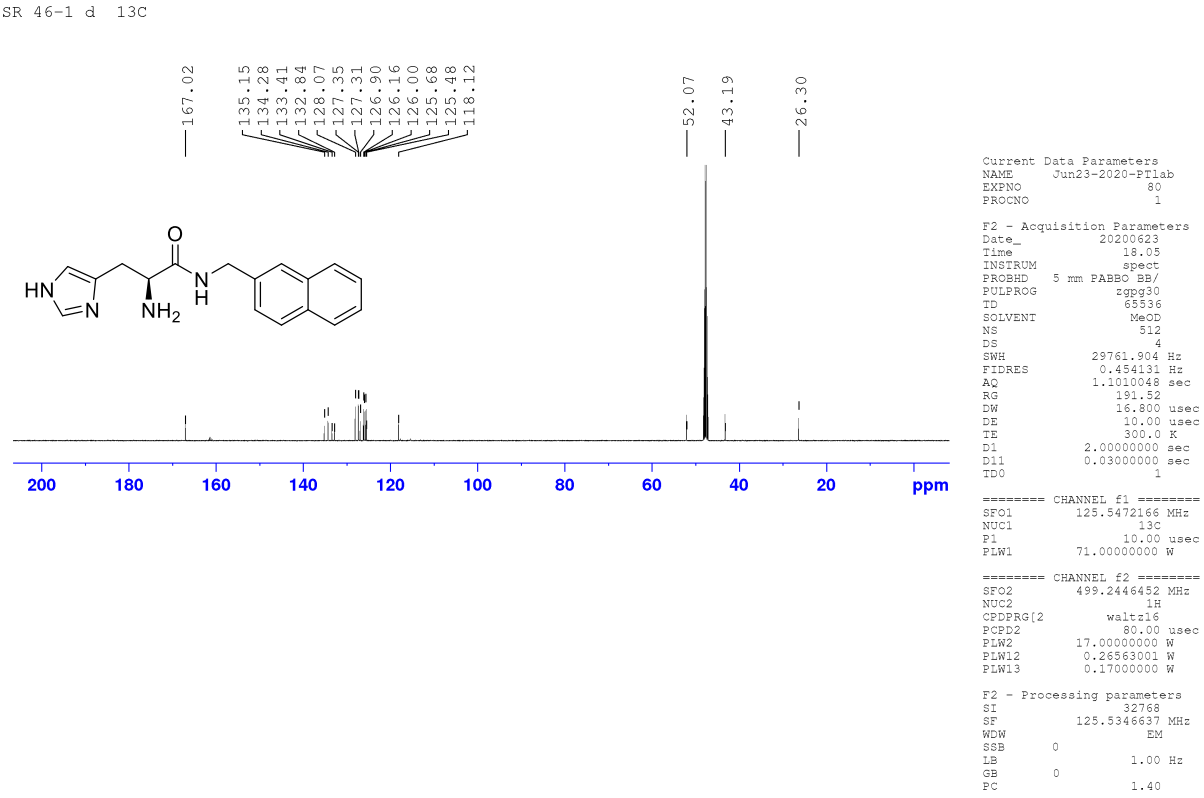
^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(naphthalen-2-yl)ethyl)propanamide **(9g):**



^{1}H NMR (500 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-2-ylmethyl)propanamide **(10a):**



^{13}C NMR (126 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-2-ylmethyl)propanamide **(10a):**




```
Current Data Parameters
NAME      Jun23-2020-PTlab
EXPNO                   80
PROCNO                   1

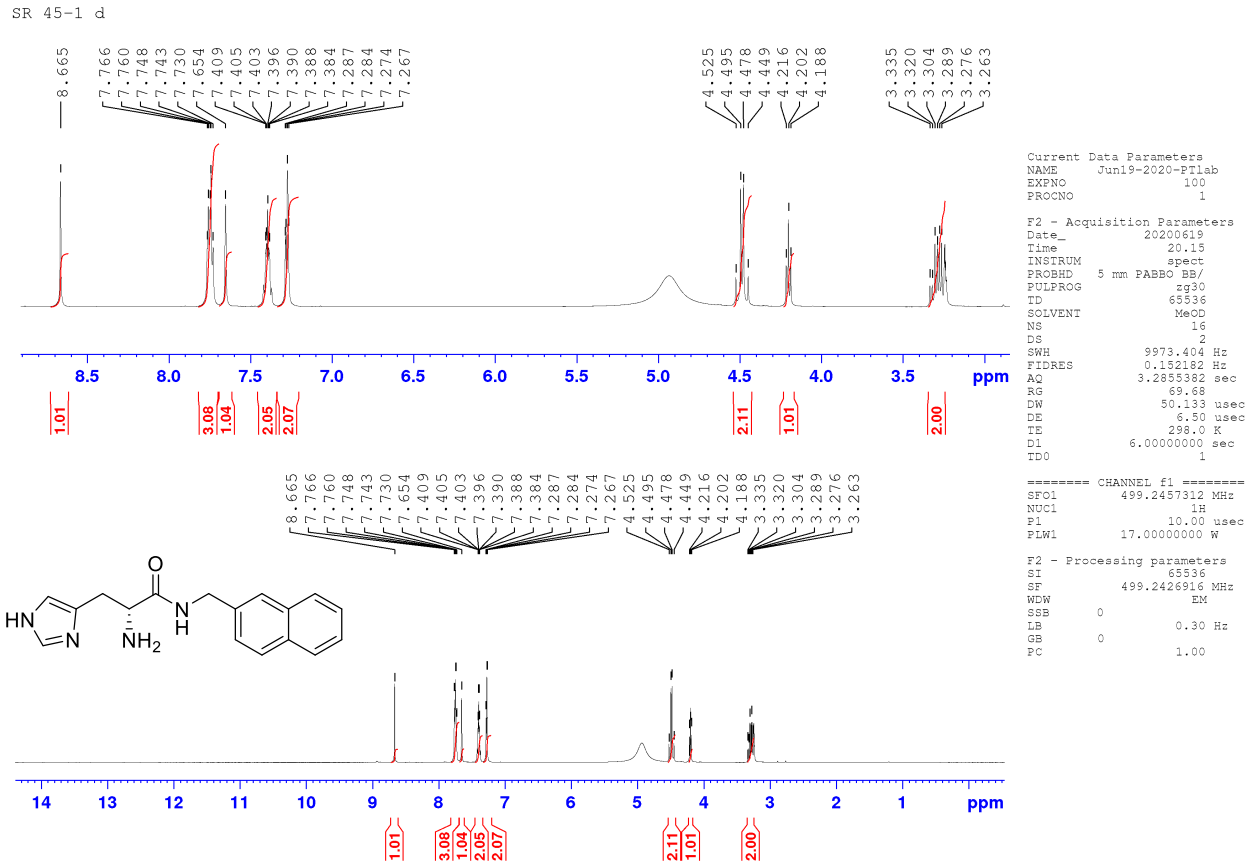
F2 - Acquisition Parameters
Date_               20200623
Time                   18.05
INSTRUM                spect
PROBHD    5 mm PABBO BB/
PULPROG             zgpg30
TD                   65536
SOLVENT               MeOD
NS                     512
DS                       4
SWH              29761.904 Hz
FIDRES           0.454131 Hz
AQ               1.1010048 sec
RG                  191.52
DW                 16.800 usec
DE                 10.00 usec
TE                 300.0 K
D1              2.00000000 sec
D11             0.03000000 sec
TD0                      1

======== CHANNEL f1 ========
SFO1          125.5472166 MHz
NUC1                   13C
P1                   10.00 usec
PLW1          71.00000000 W

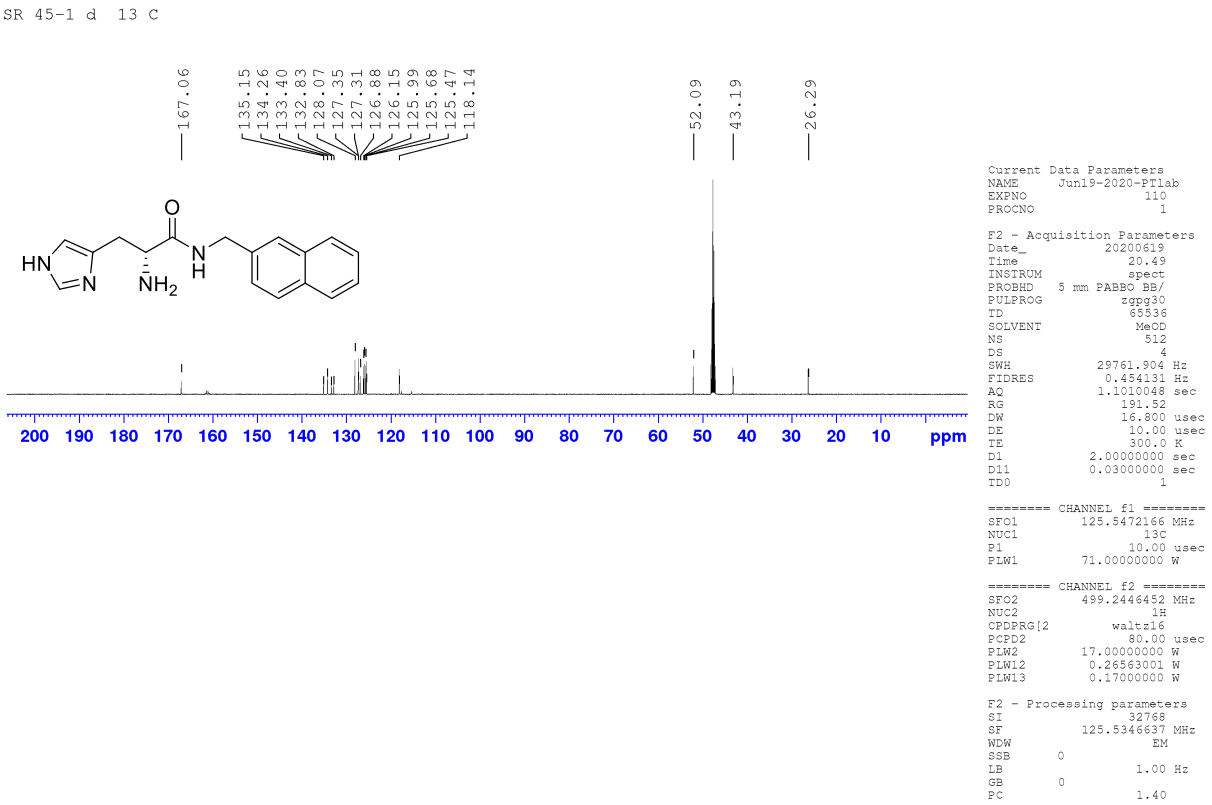
======== CHANNEL f2 ========
SFO2          499.2446452 MHz
NUC2                   1H
CPDPRG[2           waltz16
PCPD2                80.00 usec
PLW2          17.00000000 W
PLW12          0.26563001 W
PLW13          0.17000000 W

F2 - Processing parameters
SI                   32768
SF            125.5346637 MHz
WDW                     EM
SSB       0
LB                   1.00 Hz
GB        0
PC                   1.40
```

^{1}H NMR (500 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-2-ylmethyl)propanamide **(10b):**



^{13}C NMR (126 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-2-ylmethyl)propanamide **(10b):**




```
Current Data Parameters
NAME      Jun19-2020-PTlab
EXPNO                 110
PROCNO                  1

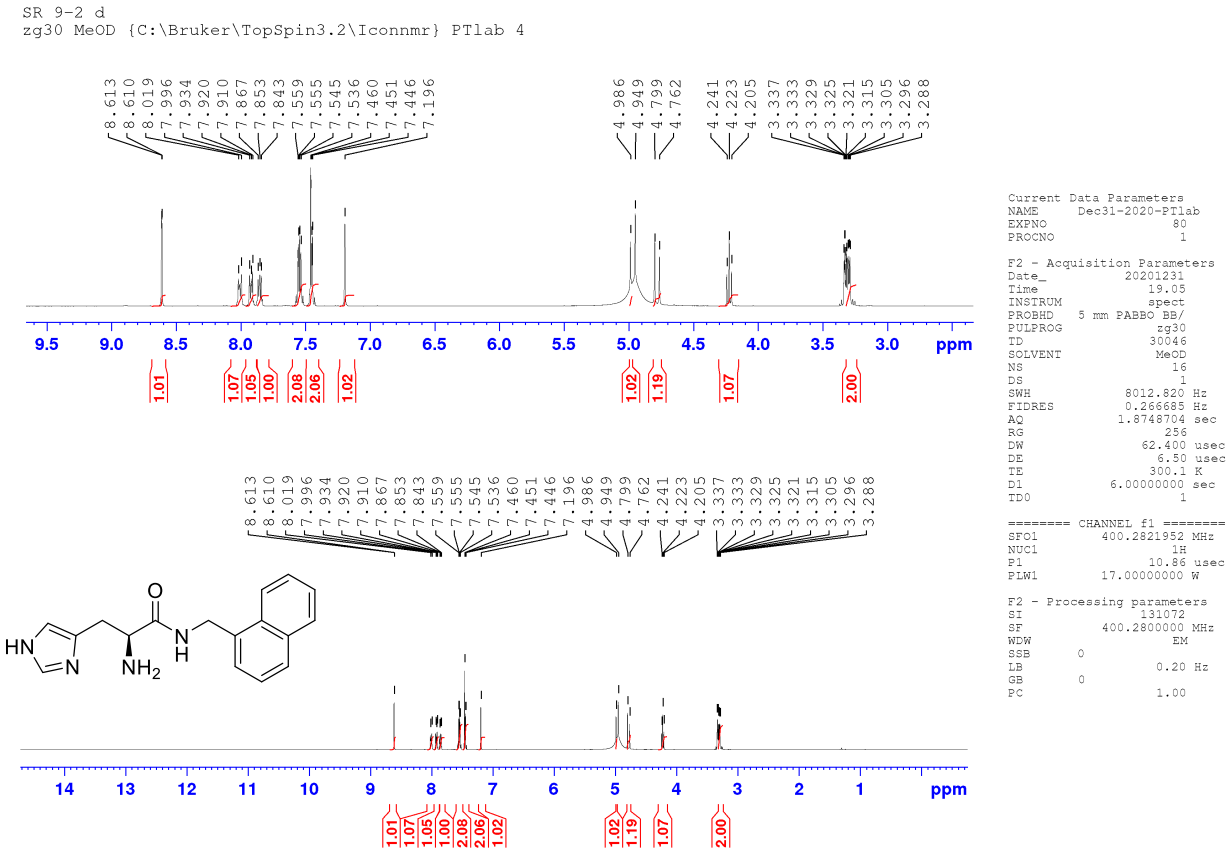
F2 - Acquisition Parameters
Date_              20200619
Time                 20.49
INSTRUM              spect
PROBHD   5 mm PABBO BB/
PULPROG            zgpg30
TD                  65536
SOLVENT              MeOD
NS                    512
DS                      4
SWH              29761.904 Hz
FIDRES            0.454131 Hz
AQ                1.1010048 sec
RG                  191.52
DW                 16.800 usec
DE                 10.00 usec
TE                 300.0 K
D1              2.00000000 sec
D11             0.03000000 sec
TD0                     1

======== CHANNEL f1 ========
SFO1          125.5472166 MHz
NUC1                    13C
P1                   10.00 usec
PLW1          71.00000000 W

======== CHANNEL f2 ========
SFO2          499.2446452 MHz
NUC2                    1H
CPDPRG[2           waltz16
PCPD2                80.00 usec
PLW2          17.00000000 W
PLW12          0.26563001 W
PLW13          0.17000000 W

F2 - Processing parameters
SI                   32768
SF            125.5346637 MHz
WDW                     EM
SSB       0
LB                    1.00 Hz
GB        0
PC                    1.40
```

^{1}H NMR (400 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-1-ylmethyl)propanamide **(10c):**



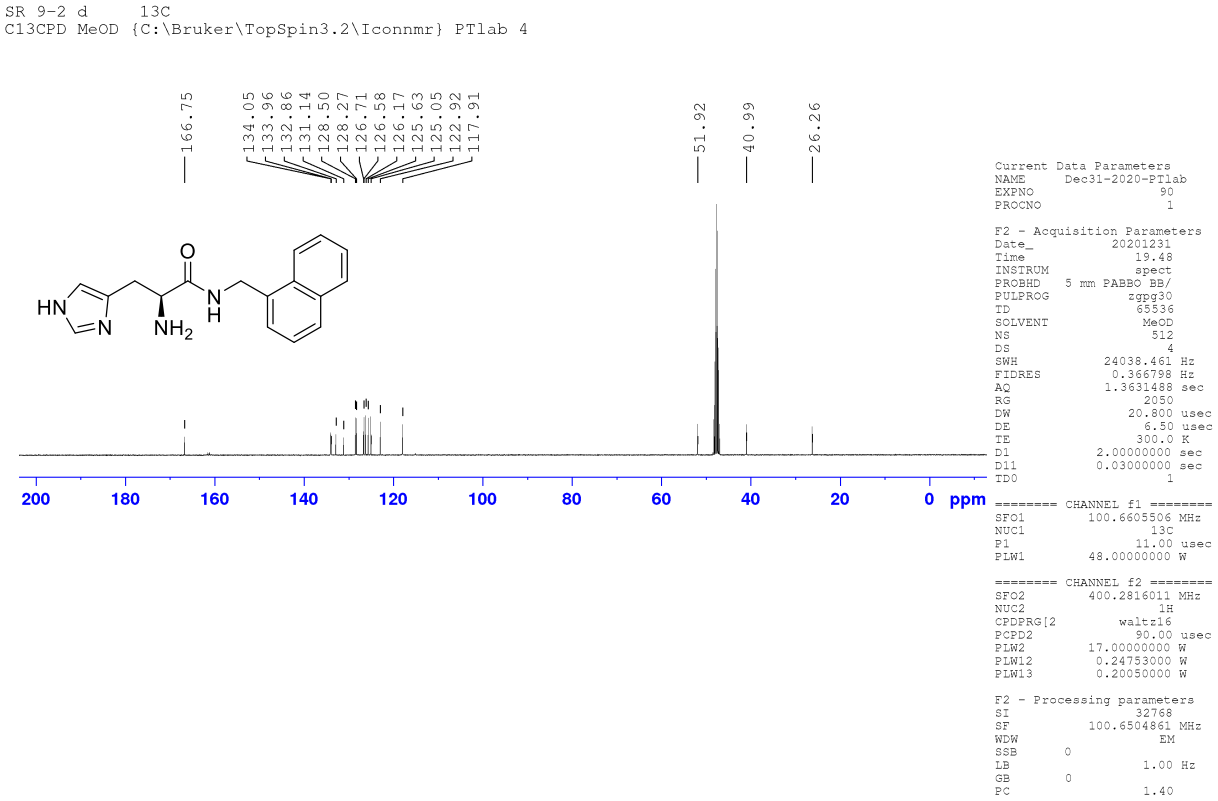

```
Current Data Parameters
NAME      Dec31-2020-PTlab
EXPNO                   80
PROCNO                   1

F2 - Acquisition Parameters
Date_              20201231
Time                 19.05
INSTRUM              spect
PROBHD    5 mm PABBO BB/
PULPROG              zg30
TD                  30046
SOLVENT              MeOD
NS                     16
DS                      1
SWH              8012.820 Hz
FIDRES           0.266685 Hz
AQ               1.8748704 sec
RG                    256
DW                 62.400 usec
DE                  6.50 usec
TE                 300.1 K
D1              6.00000000 sec
TD0                     1

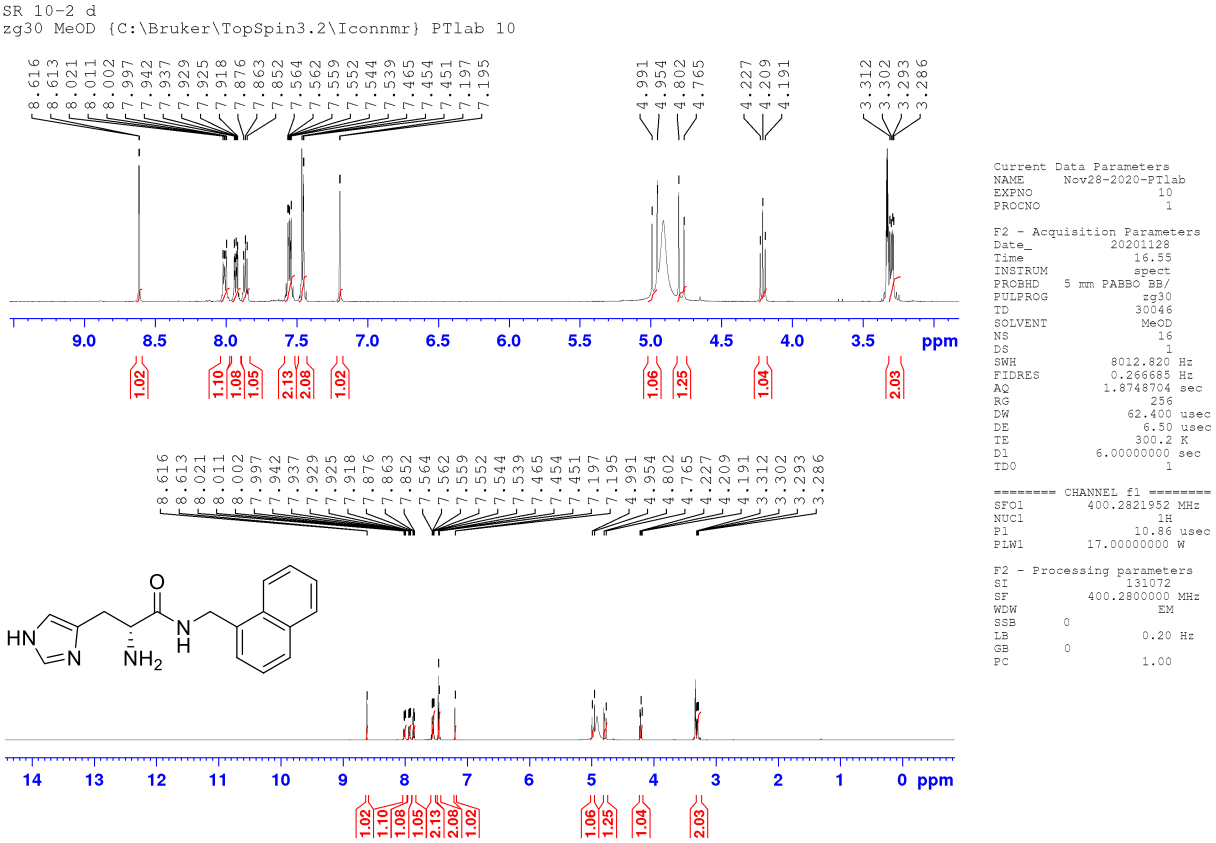
======== CHANNEL f1 ========
SFO1          400.2821952 MHz
NUC1                    1H
P1                   10.86 usec
PLW1         17.00000000 W

F2 - Processing parameters
SI                   131072
SF            400.2800000 MHz
WDW                     EM
SSB       0
LB                   0.20 Hz
GB        0
PC                   1.00
```

^{13}C NMR (100 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-1-ylmethyl)propanamide **(10c):**



^{1}H NMR (400 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-1-ylmethyl)propanamide **(10d):**



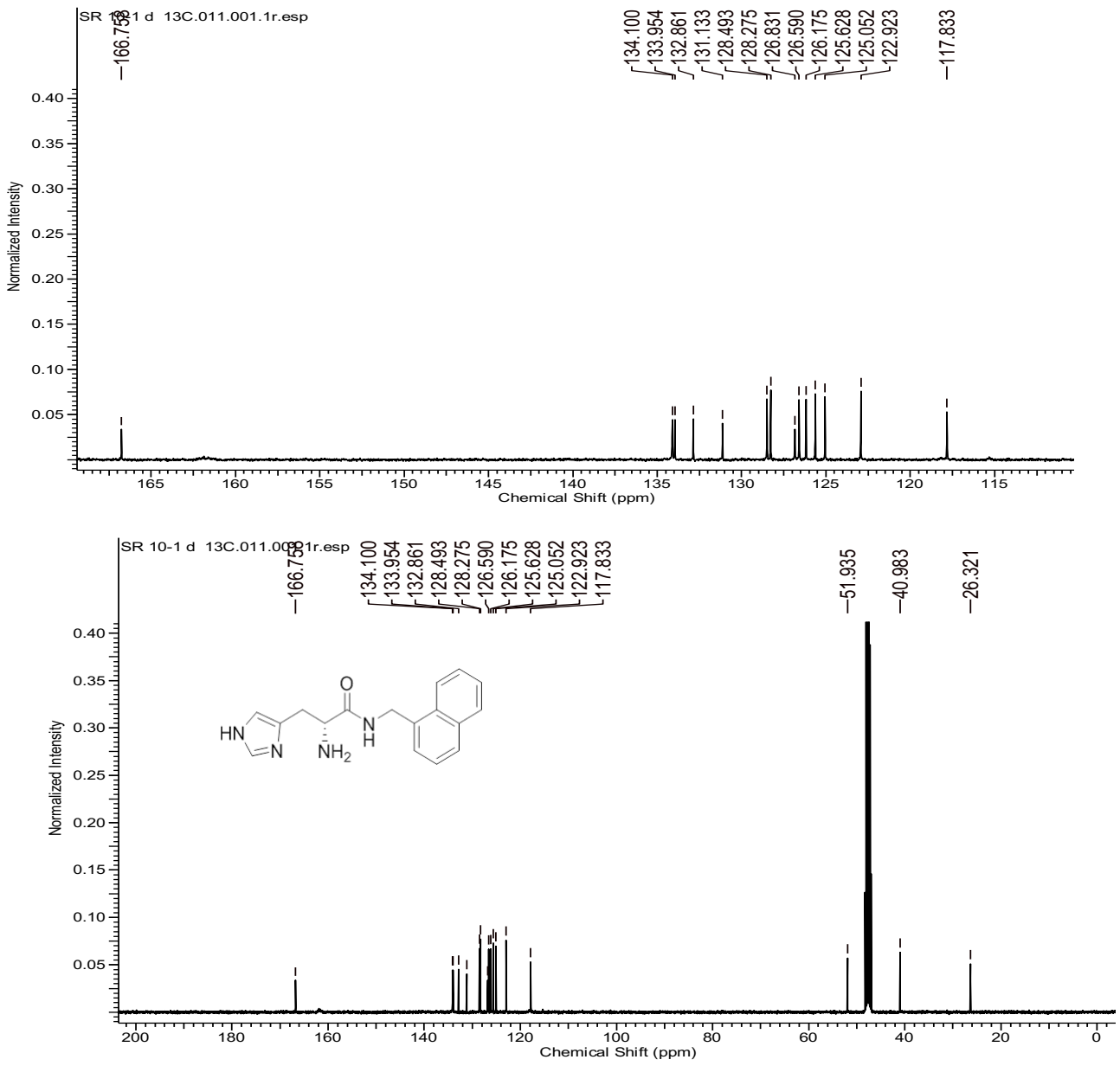

```
Current Data Parameters
NAME      Nov28-2020-PTlab
EXPNO                   10
PROCNO                   1

F2 - Acquisition Parameters
Date_              20201128
Time                 16.55
INSTRUM              spect
PROBHD   5 mm PABBO BB/
PULPROG              zg30
TD                  30046
SOLVENT              MeOD
NS                     16
DS                      1
SWH              8012.820 Hz
FIDRES           0.266685 Hz
AQ               1.8748704 sec
RG                    256
DW                 62.400 usec
DE                  6.50 usec
TE                 300.2 K
D1               6.00000000 sec
TD0                     1

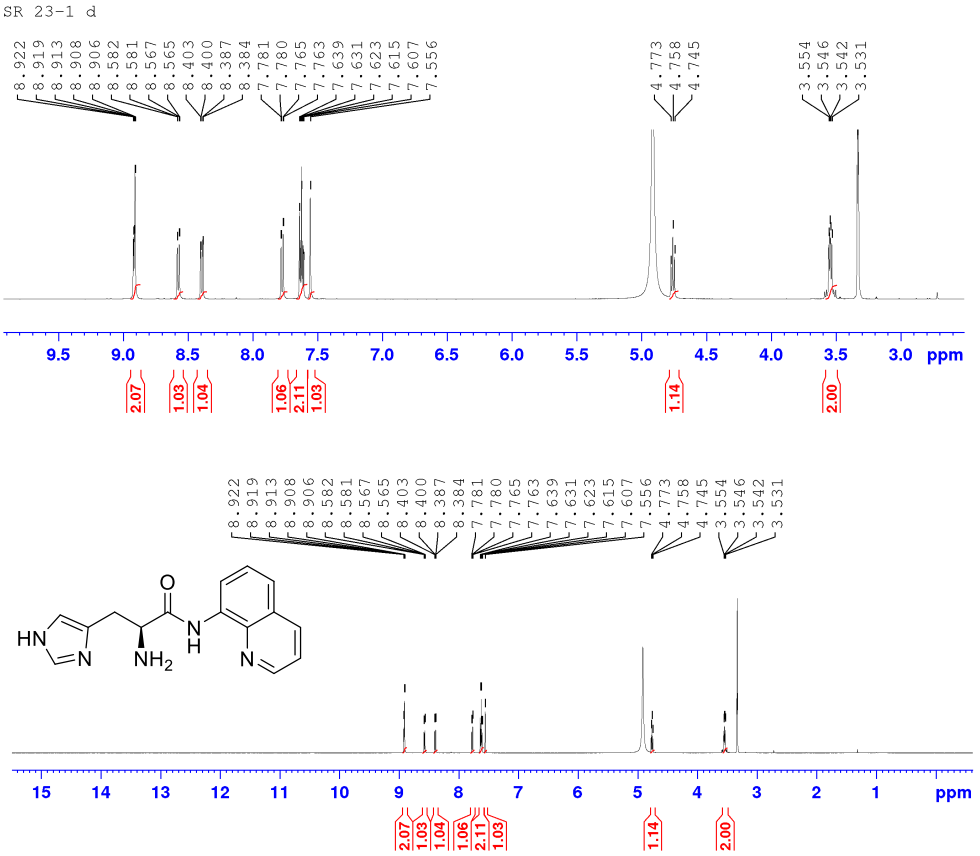
======== CHANNEL f1 ========
SFO1         400.2821952 MHz
NUC1                    1H
P1                  10.86 usec
PLW1        17.00000000 W

F2 - Processing parameters
SI                  131072
SF           400.2800000 MHz
WDW                     EM
SSB       0
LB                  0.20 Hz
GB        0
PC                  1.00
```

^{13}C NMR (100 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-1-ylmethyl)propanamide **(10d):**



^{1}H NMR (500 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(quinolin-8-yl)propanamide **(11a)*:***




```
Current Data Parameters
NAME      Nov15-2019-PTlab
EXPNO                   20
PROCNO                   1

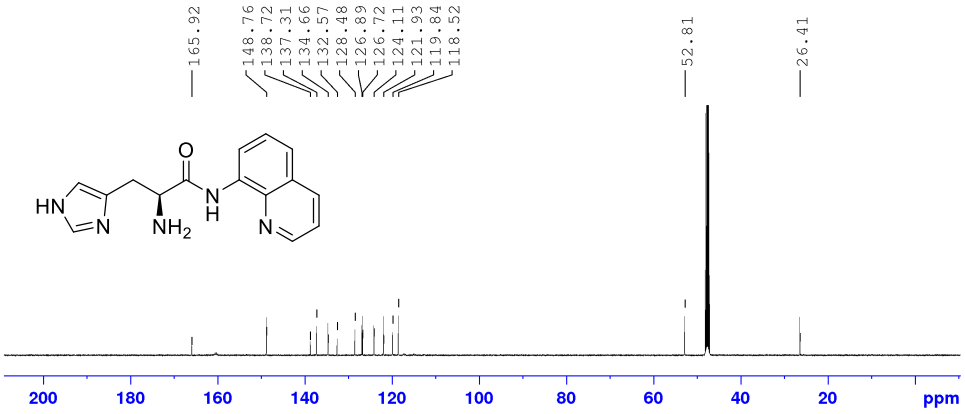
F2 - Acquisition Parameters
Date_            20191115
Time                19.21
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG             zg30
TD                 65536
SOLVENT             MeOD
NS                    16
DS                     2
SWH             9973.404 Hz
FIDRES          0.152182 Hz
AQ             3.2855382 sec
RG                152.34
DW                50.133 usec
DE                 6.50 usec
TE                298.0 K
D1             6.00000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1          499.2621482 MHz
NUC1                   1H
P1                  10.00 usec
PLW1          17.00000000 W

F2 - Processing parameters
SI                 65536
SF            499.2590651 MHz
WDW                   EM
SSB       0
LB                 0.30 Hz
GB        0
PC                 1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(quinolin-8-yl)propanamide **(11a):**

SR 23-1 d 13 C




```
Current Data Parameters
NAME      Nov15-2019-PTlab
EXPNO                   50
PROCNO                   1

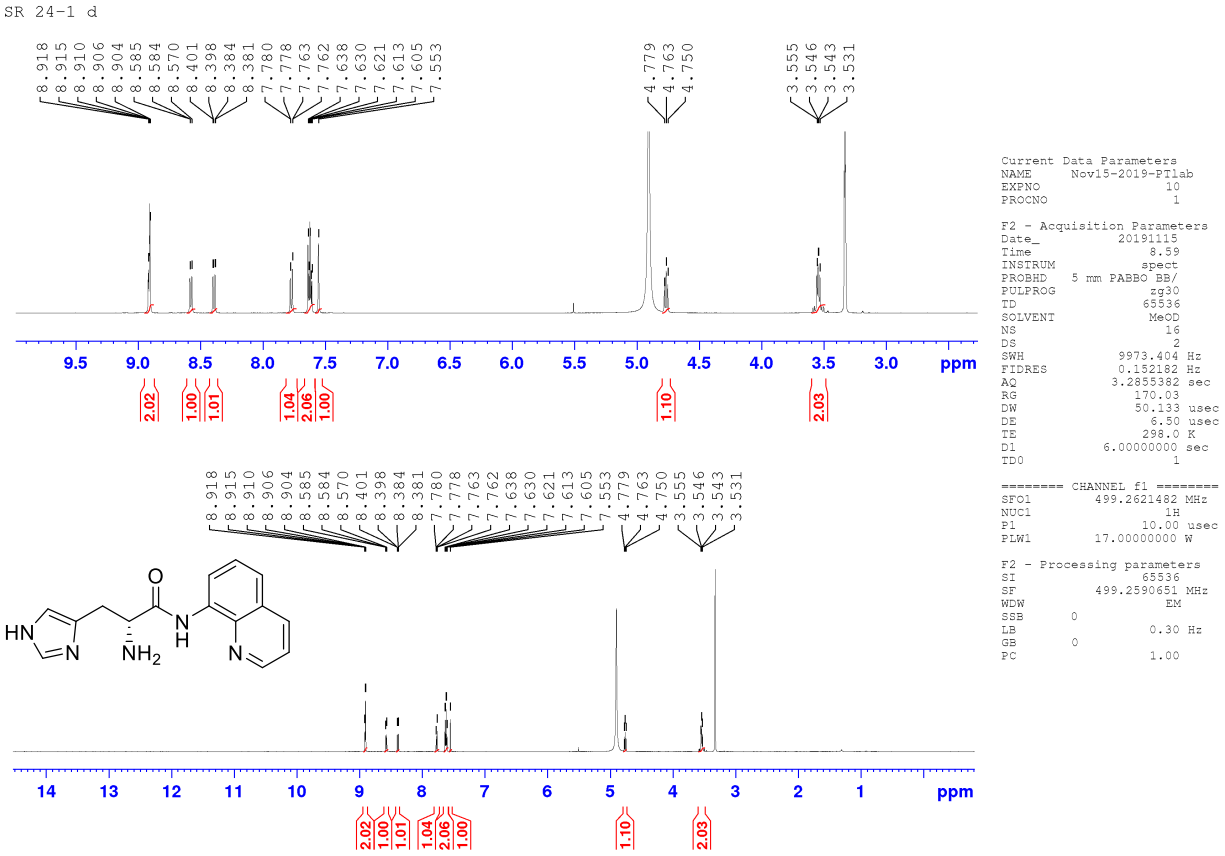
F2 - Acquisition Parameters
Date_               20191115
Time                   20.56
INSTRUM                spect
PROBHD    5 mm PABBO BB/
PULPROG             zgpg30
TD                   65536
SOLVENT               MeOD
NS                     512
DS                       4
SWH              29761.904 Hz
FIDRES            0.454131 Hz
AQ                1.1010048 sec
RG                   191.52
DW                  16.800 usec
DE                  10.00 usec
TE                  300.0 K
D1             2.00000000 sec
D11            0.03000000 sec
TD0                      1

======== CHANNEL f1 ========
SFO1          125.5513451 MHz
NUC1                    13C
P1                    10.00 usec
PLW1          71.00000000 W

======== CHANNEL f2 ========
SFO2          499.2610621 MHz
NUC2                    1H
CPDPRG[2           waltz16
PCPD2                 80.00 usec
PLW2          17.00000000 W
PLW12          0.26563001 W
PLW13          0.17000000 W

F2 - Processing parameters
SI                     32768
SF            125.5387918 MHz
WDW                       EM
SSB       0
LB                     1.00 Hz
GB        0
PC                     1.40
```

^{1}H NMR (500 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(quinolin-8-yl)propanamide **(11b):**



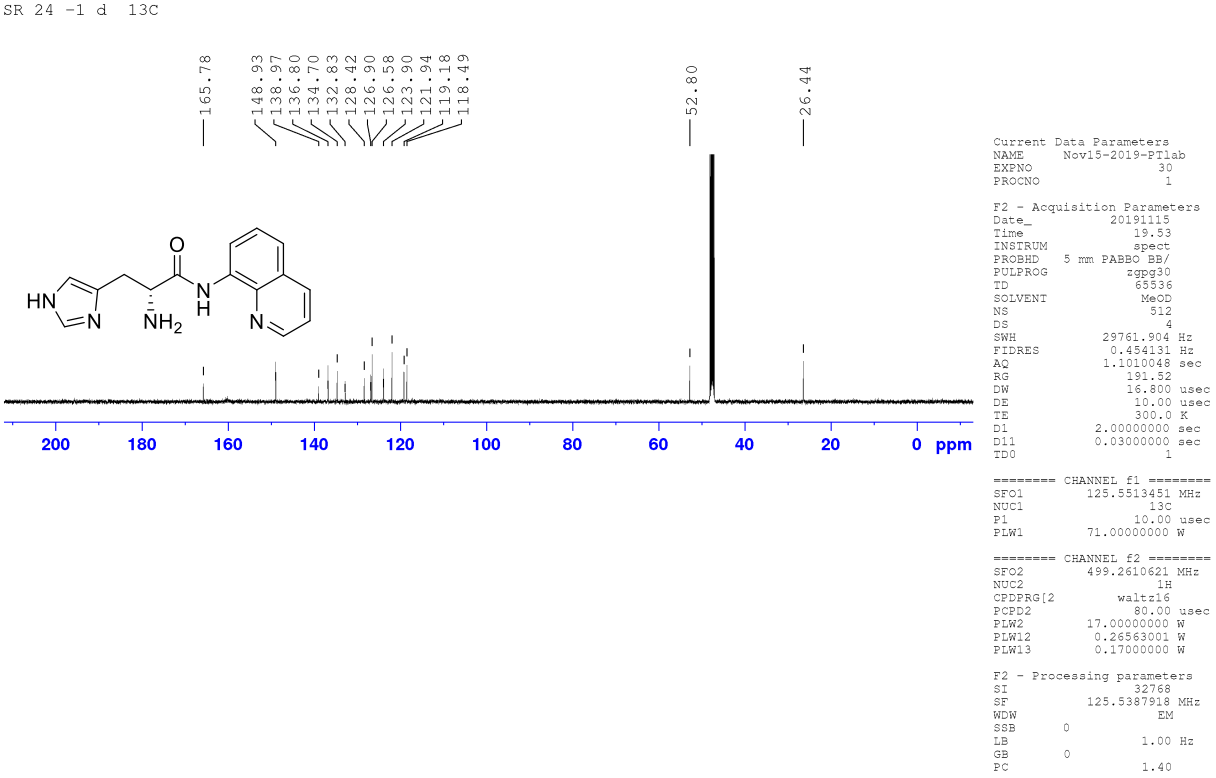

```
Current Data Parameters
NAME      Nov15-2019-PTlab
EXPNO                   10
PROCNO                   1

F2 - Acquisition Parameters
Date_            20191115
Time                 8.59
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG             zg30
TD                 65536
SOLVENT             MeOD
NS                    16
DS                     2
SWH             9973.404 Hz
FIDRES          0.152182 Hz
AQ              3.2855382 sec
RG                 170.03
DW                50.133 usec
DE                 6.50 usec
TE                298.0 K
D1             6.00000000 sec
TD0                    1

======== CHANNEL f1 ========
SFO1         499.2621482 MHz
NUC1                   1H
P1                10.00 usec
PLW1         17.00000000 W

F2 - Processing parameters
SI                 65536
SF           499.2590651 MHz
WDW                   EM
SSB       0
LB                0.30 Hz
GB        0
PC                1.00
```

^{13}C NMR (126 MHz, MeOH-d_4), (*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(quinolin-8-yl)propanamide **(11b):**




```
Current Data Parameters
NAME      Nov15-2019-PTlab
EXPNO                   30
PROCNO                   1

F2 - Acquisition Parameters
Date_            20191115
Time                19.53
INSTRUM             spect
PROBHD   5 mm PABBO BB/
PULPROG          zgpg30
TD                65536
SOLVENT             MeOD
NS                    512
DS                      4
SWH           29761.904 Hz
FIDRES         0.454131 Hz
AQ            1.1010048 sec
RG                191.52
DW                16.800 usec
DE                10.00 usec
TE                300.0 K
D1           2.00000000 sec
D11          0.03000000 sec
TD0                     1

======== CHANNEL f1 ========
SFO1        125.5513451 MHz
NUC1                13C
P1                10.00 usec
PLW1        71.00000000 W

======== CHANNEL f2 ========
SFO2        499.2610621 MHz
NUC2                1H
CPDPRG[2         waltz16
PCPD2             80.00 usec
PLW2        17.00000000 W
PLW12        0.26563001 W
PLW13        0.17000000 W

F2 - Processing parameters
SI                  32768
SF          125.5387918 MHz
WDW                    EM
SSB       0
LB                 1.00 Hz
GB        0
PC                 1.40
```

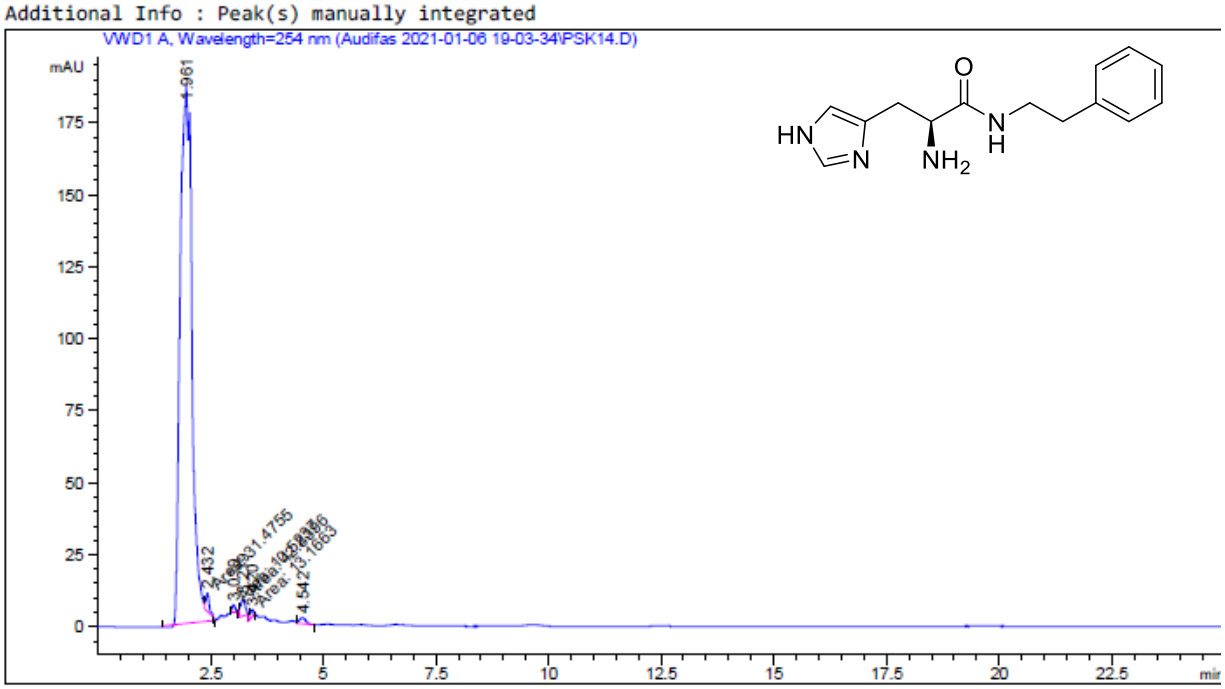
HPLC Chromatograms of Final Compounds

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenethylpropanamide **(4c)**

```
Sample Name: PSK14
======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :  28
Acq. Instrument : 1220 HPLC                       Location :  80
Injection Date  : 1/7/2021 7:33:06 AM                 Inj :  1
                                                  Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : PSK14

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.961 MM R  0.3806 3403.53442  186.91812  96.6275
   2   2.432 MM T  0.0827   31.47550    6.34564   0.8936
   3   3.010 MM T  0.0727   10.52366    2.41180   0.2988
   4   3.219 MM T  0.1242   42.83961    5.74662   1.2162
```

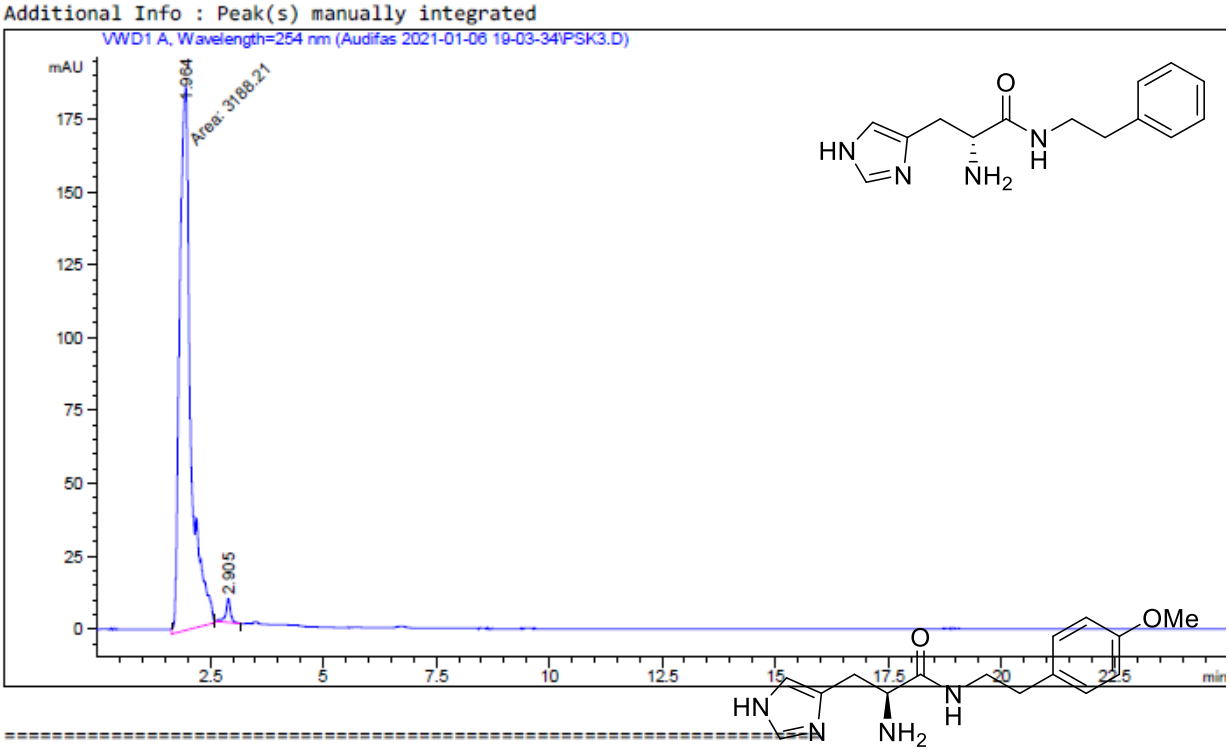
(R)-2-amino-3-(1H-imidazol-4-yl)-N-phenethylpropanamide(4d)

Sample Name: PSK

```
==================================================================
Acq. Operator   : SYSTEM                        Seq. Line :  31
Acq. Instrument : 1220 HPLC                      Location :  81
Injection Date  : 1/7/2021 8:55:48 AM                Inj :  1
                                                Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : PSK3

Additional Info : Peak(s) manually integrated
```




```
==================================================================
                      Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.964 MM T  0.3095 3188.21191  186.93040  98.1449
   2   2.905 BB    0.1058   60.26109    8.05252   1.8551

Totals :                  3248.47301  194.98293
==================================================================
                      Area Percent Report
==================================================================

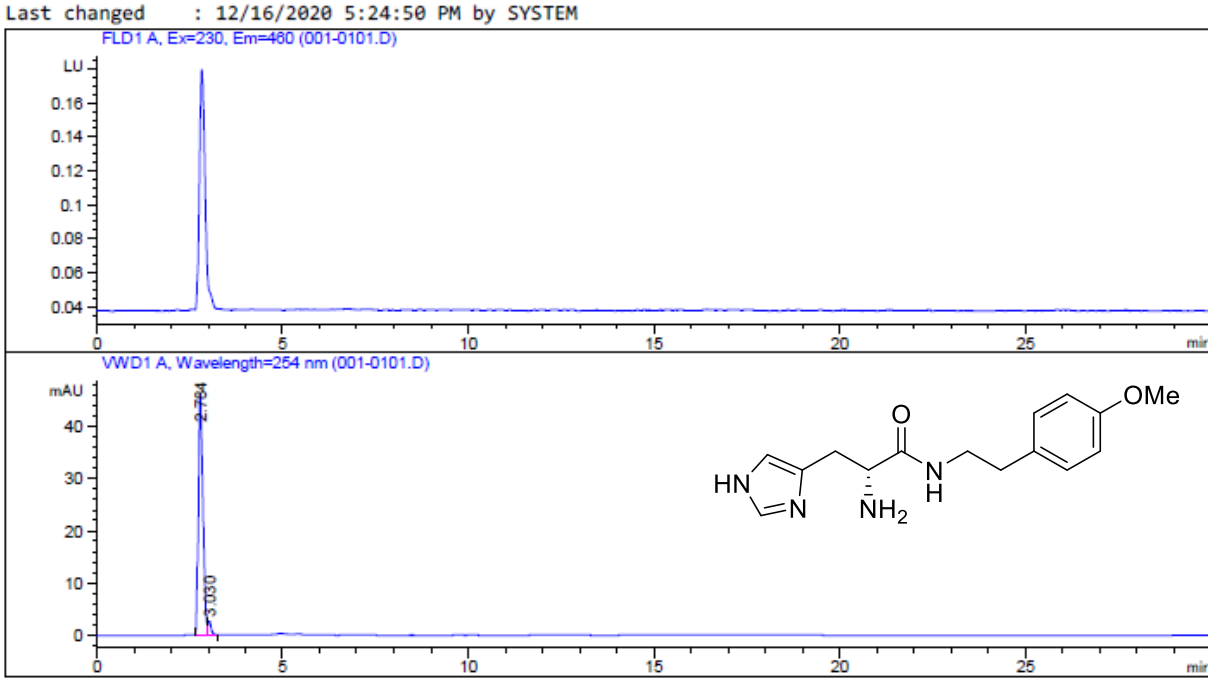
Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm
```

(R)-2-amino-3-(1H-imidazol-4-yl)-N-(4-methoxyphenethyl)propanamide (4f):

Sample Name: SR 69-1 d

```
==================================================================
  Acq. Operator   : SYSTEM                        Seq. Line :   1
  Acq. Instrument : HPLC                           Location : Vial 1
  Injection Date  : 12/16/2020 5:26:00 PM              Inj :   1
                                                  Inj Volume : 20.000 µl
  Sequence File   : C:\Chem32\1\DATA\Ao\KK 2020-12-16 17-24-50\KK.S
  Method          : C:\CHEM32\1\DATA\AO\KK 2020-12-16 17-24-50\CQ.M (Sequence Method)
  Last changed    : 12/16/2020 5:24:50 PM by SYSTEM
```




```
  ==================================================================
                          Area Percent Report
  ==================================================================

  Sorted By             :      Signal
  Multiplier            :      1.0000
  Dilution              :      1.0000
  Do not use Multiplier & Dilution Factor with ISTDs


  Signal 1: FLD1 A, Ex=230, Em=460


  Signal 2: VWD1 A, Wavelength=254 nm

  Peak RetTime Type  Width     Area      Height     Area
    #   [min]        [min]   [mAU*s]     [mAU]       %
  ----|-------|----|-------|----------|----------|--------|
     1   2.784 BV    0.1141  388.43942   46.19957  96.1296
     2   3.030 VB    0.0900   15.63965    2.60797   3.8704

  Totals :                   404.07908   48.80754


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.741 BB    0.2070 2580.59741  184.69226 100.0000

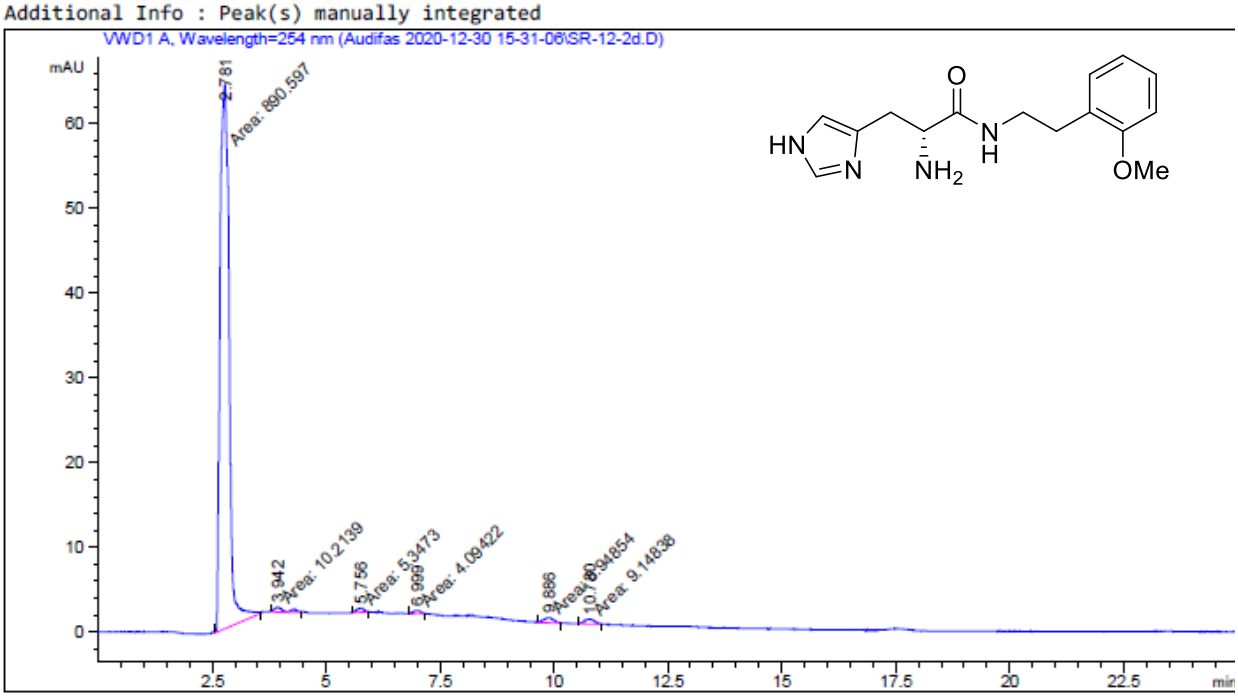
Totals :                  2580.59741  184.69226
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-methoxyphenethyl)propanamide **(4h)**

```
Sample Name: SR-12-2d
======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   1
Acq. Instrument : 1220 HPLC                       Location :   51
Injection Date  : 12/30/2020 3:32:53 PM               Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-30 15-31-06\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/30/2020 3:31:09 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05% TFA / Methanol

Sample Info     : SR-12-2d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|---------|----------|--------|
   1   2.781 MM T  0.2314  890.59698   64.13306  95.9334
   2   3.942 MM    0.2926   10.21391 5.81875e-1   1.1002
   3   5.756 MM    0.1739    5.34730 5.12450e-1   0.5760
   4   6.999 MM    0.1874    4.09422 3.64115e-1   0.4410
```

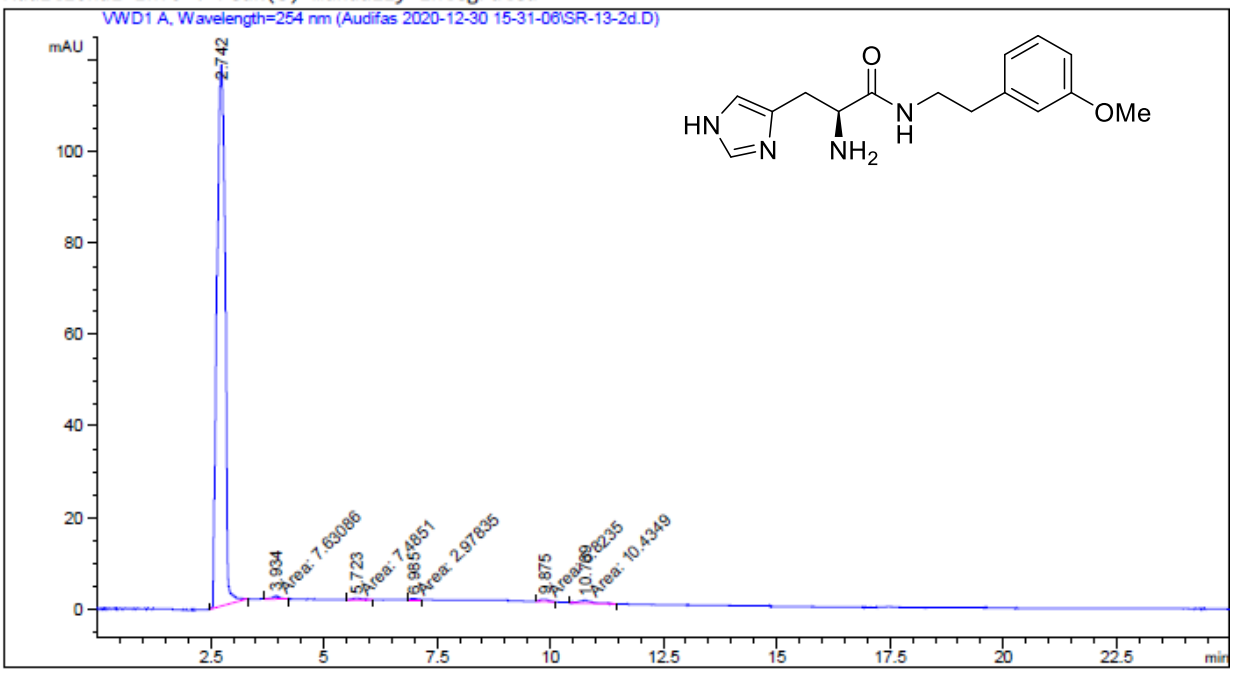
(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-methoxyphenethyl)propanamide **(4i)**

```
Sample Name: SR-13-2d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   2
Acq. Instrument : 1220 HPLC                       Location :   52
Injection Date  : 12/30/2020 4:00:19 PM              Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-30 15-31-06\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/30/2020 3:31:09 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05% TFA / Methanol

Sample Info     : SR-13-2d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs

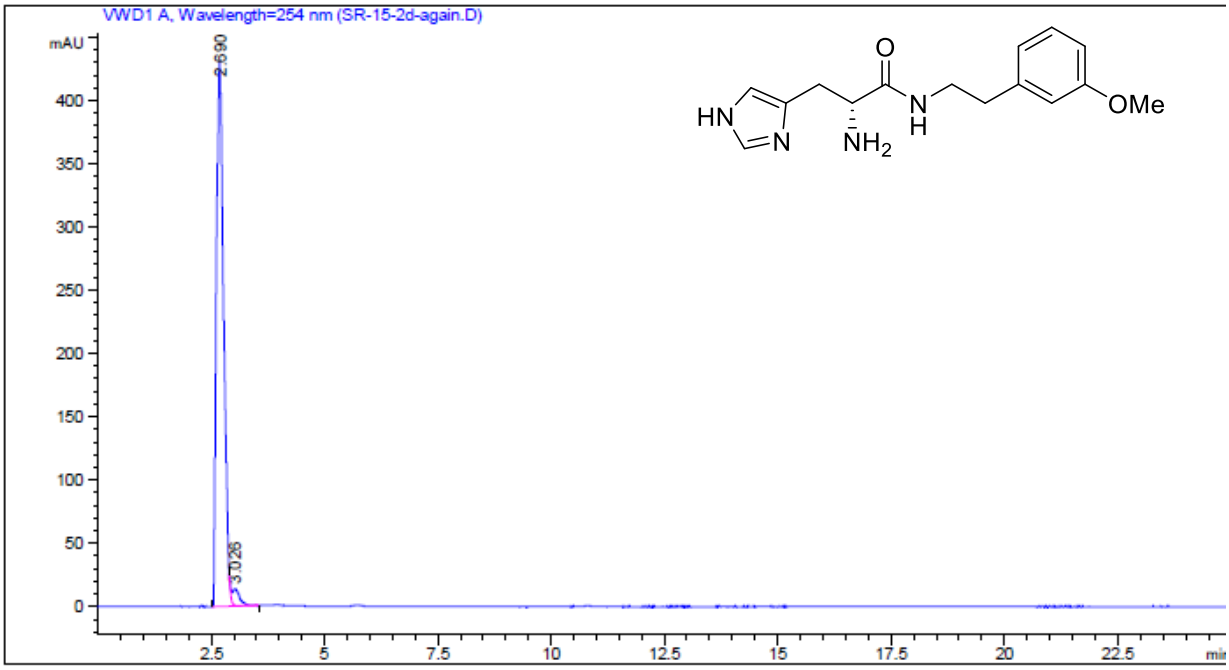

Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.742 BB    0.1802 1465.06689  118.49198  97.7089
   2   3.934 MM T  0.2183    7.63086 6.10415e-1   0.5089
   3   5.723 MM T  0.2582    7.48510 4.83206e-1   0.4992
   4   6.985 MM T  0.1941    2.97835 2.55785e-1   0.1986
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-methoxyphenethyl)propanamide **(4j)**

```
Sample Name: SR-15-2d
======================================================================
Acq. Operator   : SYSTEM                        Seq. Line :   3
Acq. Instrument : 1220 HPLC                     Location :   51
Injection Date  : 12/30/2020 2:27:07 PM             Inj :   1
                                               Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Sequence File   : C:\Chem32\1\Data\Audifas 2020-12-30 13-30-17\Audifas.S
Method          : C:\Chem32\1\Data\Audifas 2020-12-30 13-30-17\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/30/2020 1:30:21 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.1 %TFA 0.05% TFA/ Methanol

Sample Info     : SR-15-2d
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.690 BV R  0.1534 4793.86426  430.65921  96.8621
   2   3.026 VB E  0.1489  155.30142   13.38018   3.1379

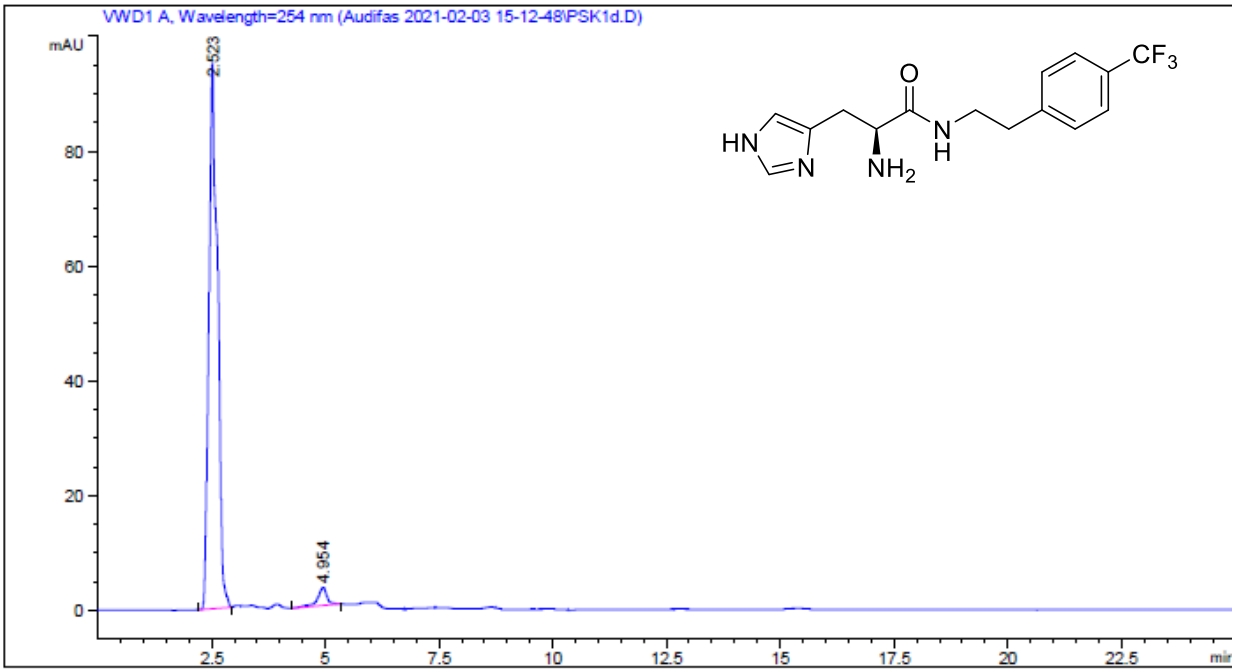
Totals :                  4949.16568  444.03939
```

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)phenethyl)propenamide

```
Sample Name: PSK1d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   8
Acq. Instrument : 1220 HPLC                       Location :   57
Injection Date  : 2/3/2021 6:26:56 PM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-02-03 15-12-48\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 2/3/2021 3:12:52 PM by SYSTEM
Method Info     : Methanol/Water 0.05 % TFA

Sample Info     : PSK1d
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs

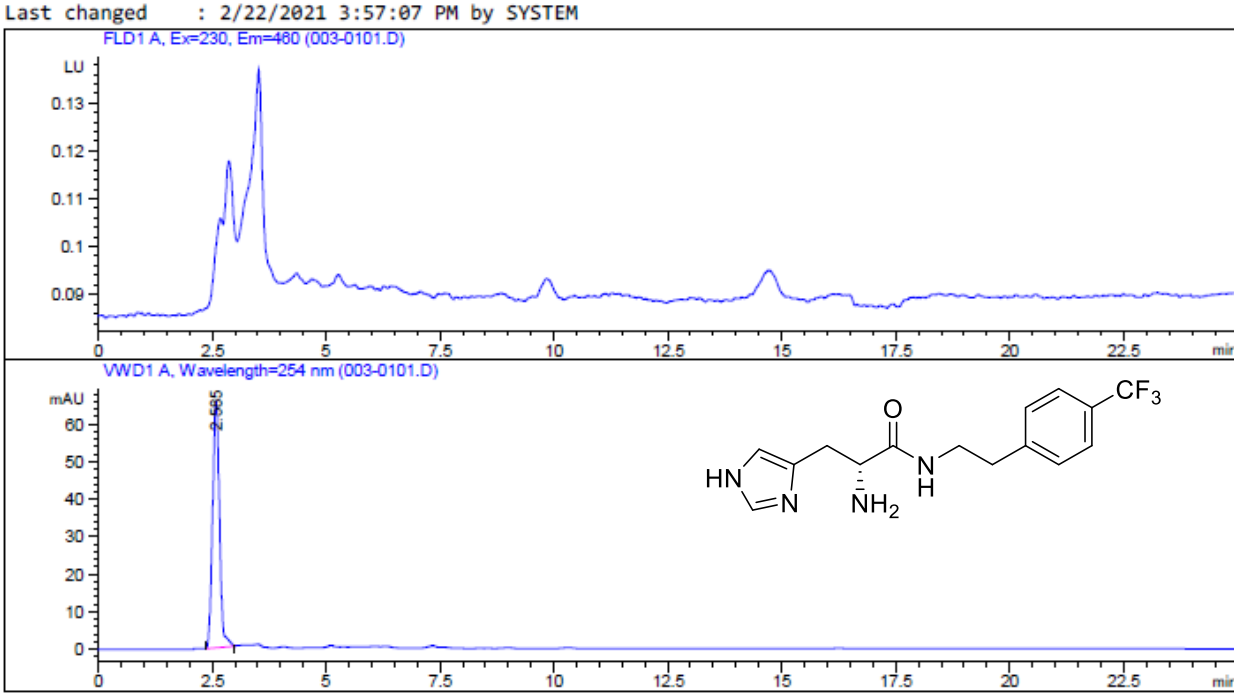

Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.523 BB    0.1667 1255.20386   95.25862  96.4839
   2   4.954 BV R  0.1736   45.74269    3.13797   3.5161

Totals :                  1300.94654   98.39659
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)phenethyl)propanamide

```
Sample Name: PSK 2d
======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   1
Acq. Instrument : HPLC                             Location : Vial 3
Injection Date  : 2/22/2021 3:58:14 PM                 Inj :   1
                                                 Inj Volume : 20.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2021-02-22 15-57-07\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2021-02-22 15-57-07\CQ.M (Sequence Method)
Last changed    : 2/22/2021 3:57:07 PM by SYSTEM
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=230, Em=460


Signal 2: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.585 BB    0.1574  694.42896   65.67966 100.0000

Totals :                   694.42896   65.67966
```

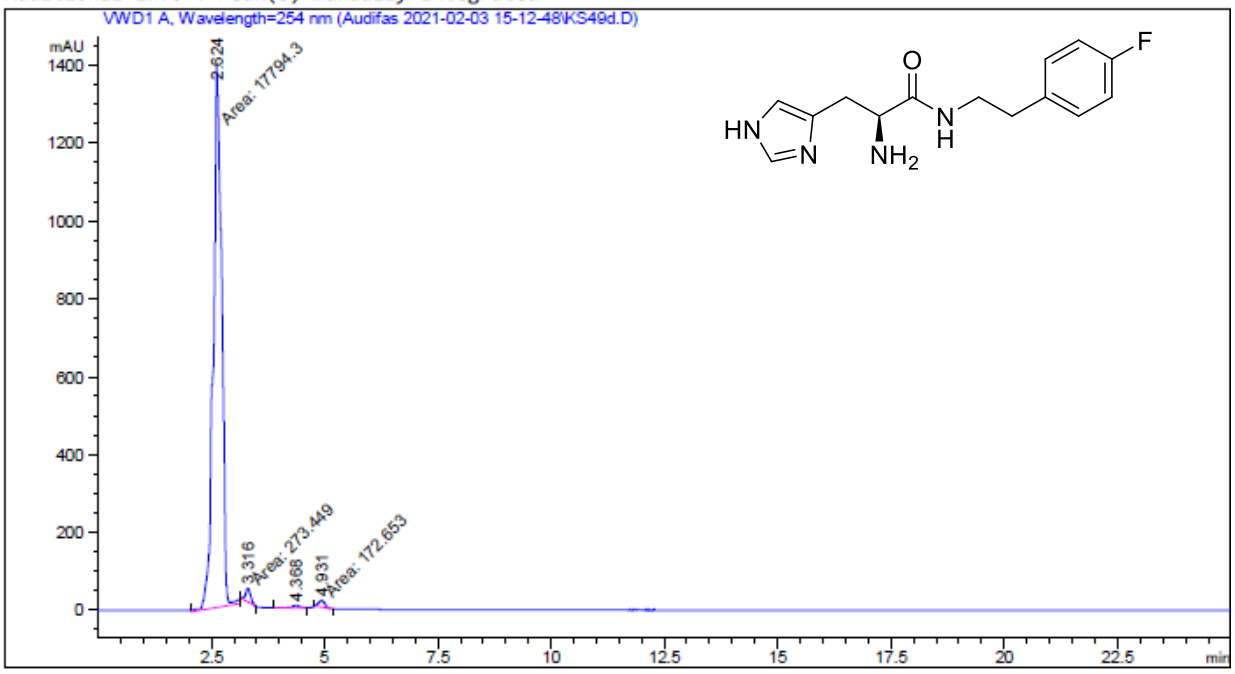
(*S*)-2-amino-*N*-(4-fluorophenethyl)-3-(1*H*-imidazol-4-yl)propanamide **(4m)**

```
Sample Name: KS49d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   6
Acq. Instrument : 1220 HPLC                       Location :   55
Injection Date  : 2/3/2021 5:31:52 PM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-02-03 15-12-48\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 2/3/2021 3:12:52 PM by SYSTEM
Method Info     : Methanol/Water 0.05 % TFA

Sample Info     : KS49d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs

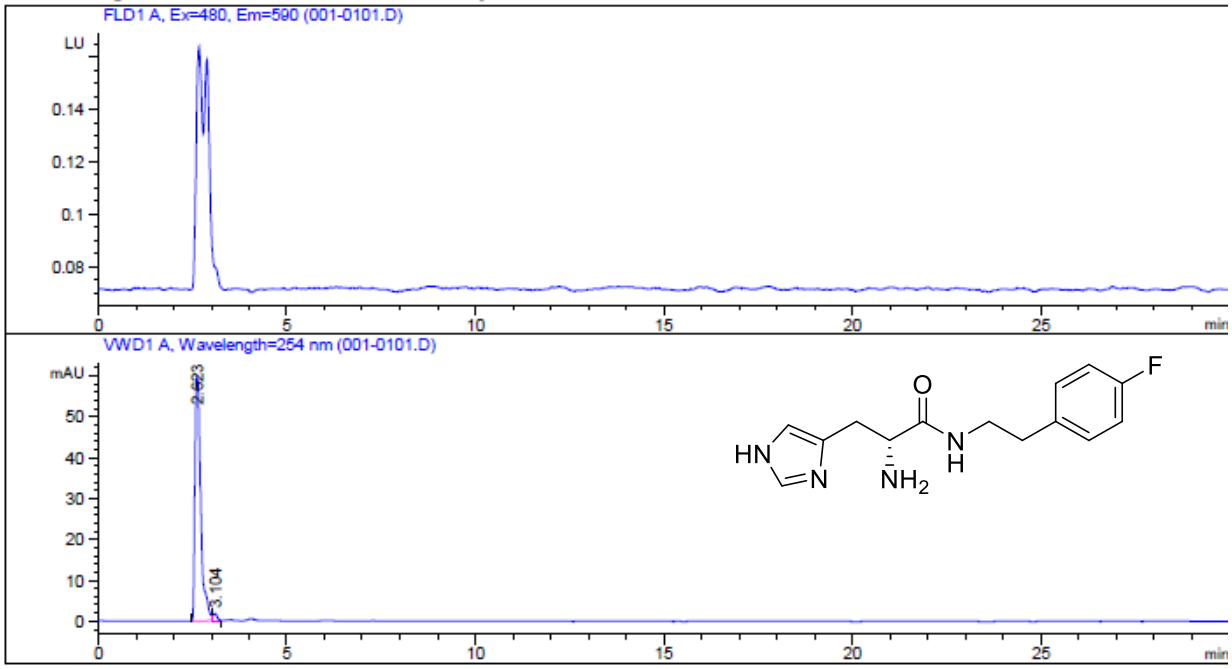

Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.624 MM T  0.2126 1.77943e4  1394.69470  96.9523
   2   3.316 MM T  0.1296  273.44873   35.17678   1.4899
   3   4.368 VV R  0.2045  113.26699    6.54201   0.6171
   4   4.931 MM T  0.1663  172.65331   17.30216   0.9407
```

(*R*)-2-amino-*N*-(4-fluorophenethyl)-3-(1*H*-imidazol-4-yl)propanamide **(4n)**

```
Sample Name: KS-94-d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   1
Acq. Instrument : HPLC                           Location : Vial 1
Injection Date  : 12/9/2020 1:13:27 PM                Inj :   1
                                                 Inj Volume : 20.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2020-12-09 13-12-18\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2020-12-09 13-12-18\CQ.M (Sequence Method)
Last changed    : 12/9/2020 1:12:19 PM by SYSTEM
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=480, Em=590


Signal 2: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.623 BV    0.1662  643.78772   59.96936  97.8111
   2   3.104 VB    0.1187   14.40736    1.70374   2.1889

Totals :                   658.19508   61.67310
```

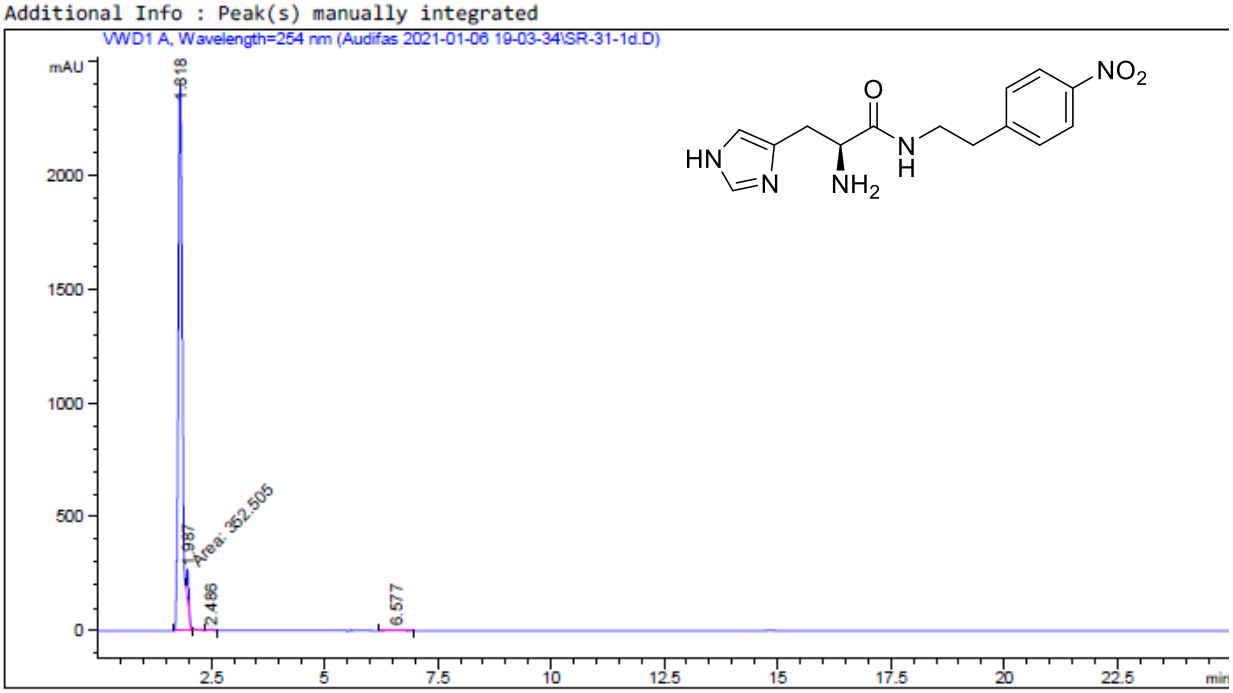
(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-nitrophenethyl)propanamide **(4o)**

```
Sample Name: SR-31-1d

=====================================================================
Acq. Operator   : SYSTEM                          Seq. Line :  13
Acq. Instrument : 1220 HPLC                        Location :  69
Injection Date  : 1/7/2021 12:39:41 AM                Inj :  1
                                                  Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR-31-1d

Additional Info : Peak(s) manually integrated
```




```
=====================================================================
                        Area Percent Report
=====================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

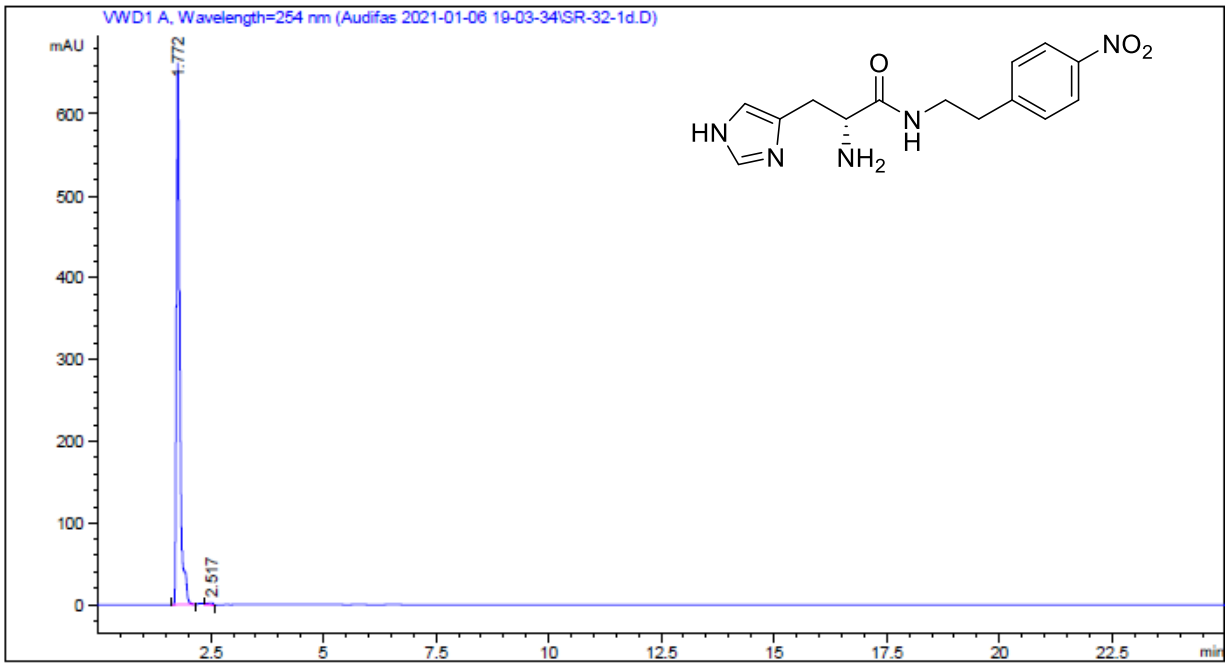
Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.818 BV R  0.0940 1.55225e4  2396.34814  97.4208
   2   1.987 MM T  0.0420  352.50507  139.91641   2.2124
   3   2.486 VB E  0.1273   33.96302    3.16218   0.2132
   4   6.577 BB    0.1635   24.48011    1.76811   0.1536
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-nitrophenethyl)propanamide **(4p)**

```
Sample Name: SR-32-1d

======================================================================
Acq. Operator   : SYSTEM                       Seq. Line :  17
Acq. Instrument : 1220 HPLC                    Location :   71
Injection Date  : 1/7/2021 2:29:57 AM               Inj :   1
                                              Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR-32-1d
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]       [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.772 BB    0.0818 3541.53955  662.76624  99.4130
   2   2.517 VB    0.1131   20.91246    2.46200   0.5870

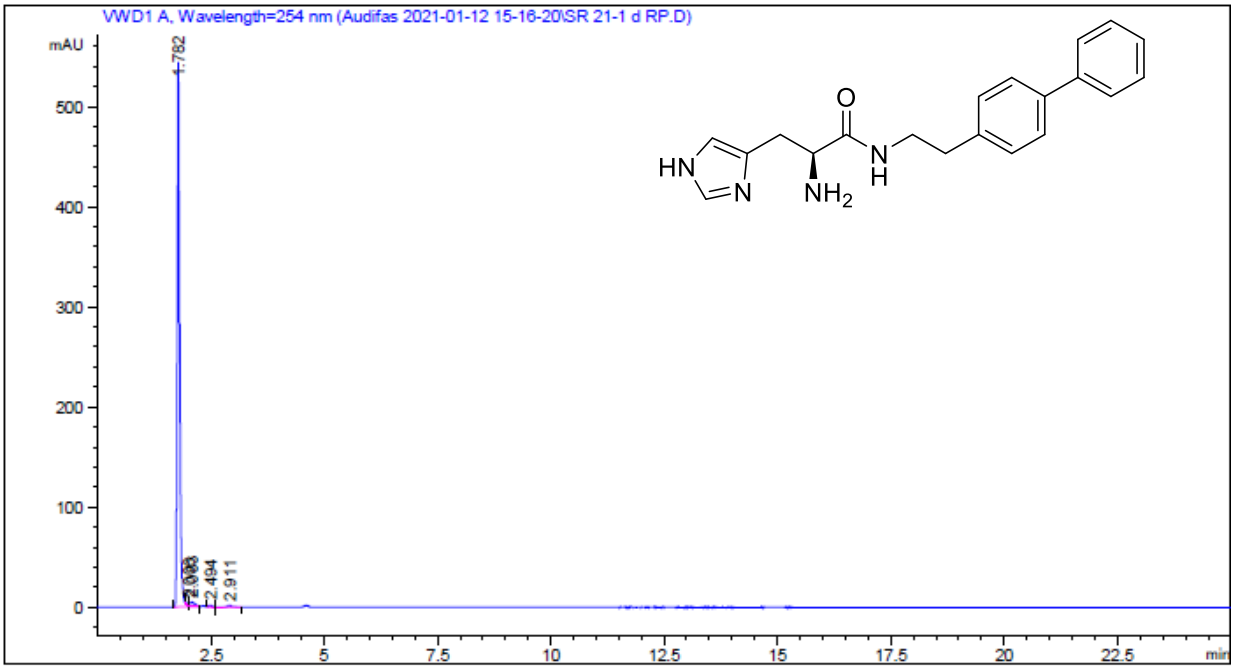
Totals :                  3562.45201  665.22823
```

(*S*)-*N*-(2-([1,1'-biphenyl]-4-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide

Sample Name: SR 21-1 d

```
==================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   4
Acq. Instrument : 1220 HPLC                     Location :   22
Injection Date  : 1/12/2021 4:30:41 PM              Inj :   1
                                               Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 10.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-12 15-16-20\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/12/2021 3:16:24 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR 21-1 d AG
```




```
==================================================================
                        Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.782 BV R  0.0580 2096.91895  544.23639  96.8984
   2   2.006 VV E  0.0471   10.06466    2.57519   0.4651
   3   2.083 VB E  0.0908   29.99245    4.43780   1.3859
   4   2.494 VB    0.0694    9.51005    1.88400   0.4395
   5   2.911 BB    0.1261   17.55167    1.74936   0.8111
```

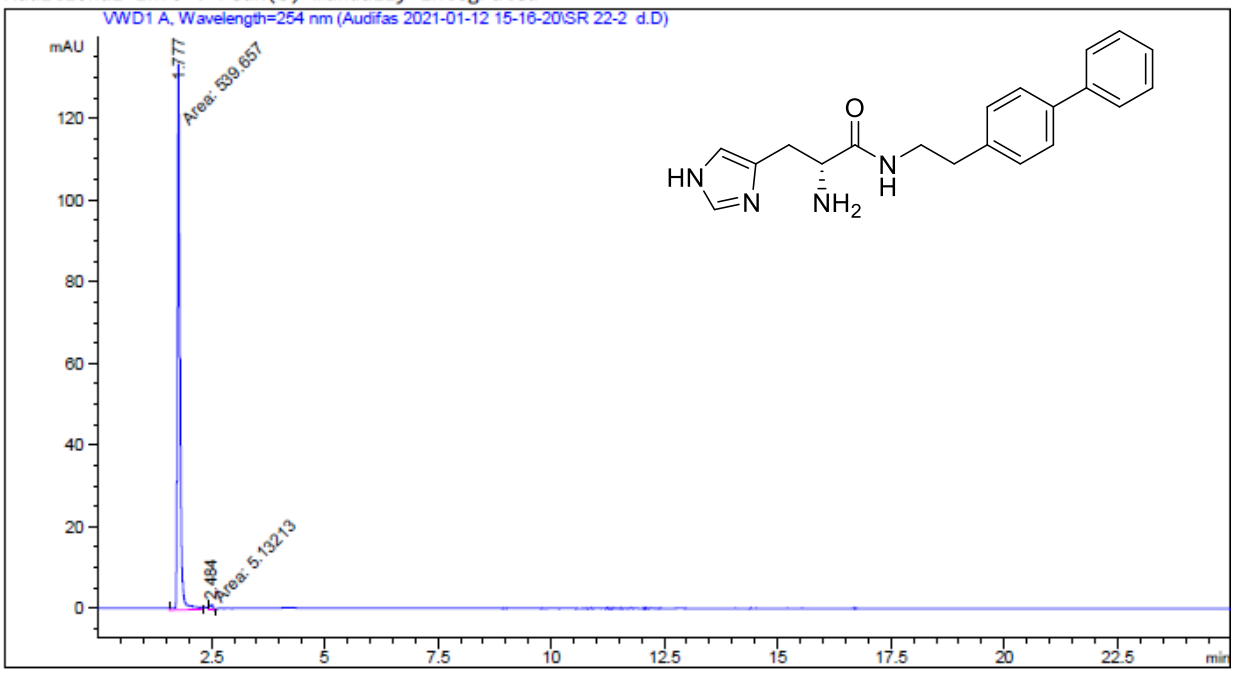
(*R*)-*N*-(2-([1,1'-biphenyl]-4-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide

Sample Name: SR 22-2 d

```
======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   5
Acq. Instrument : 1220 HPLC                       Location :   31
Injection Date  : 1/12/2021 4:57:57 PM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 10.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-12 15-16-20\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/12/2021 3:16:24 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR 22-2 d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.777 MM T  0.0673  539.65656  133.61414  99.0580
   2   2.484 MM T  0.0795    5.13213    1.07656   0.9420

Totals :                   544.78869  134.69069
```

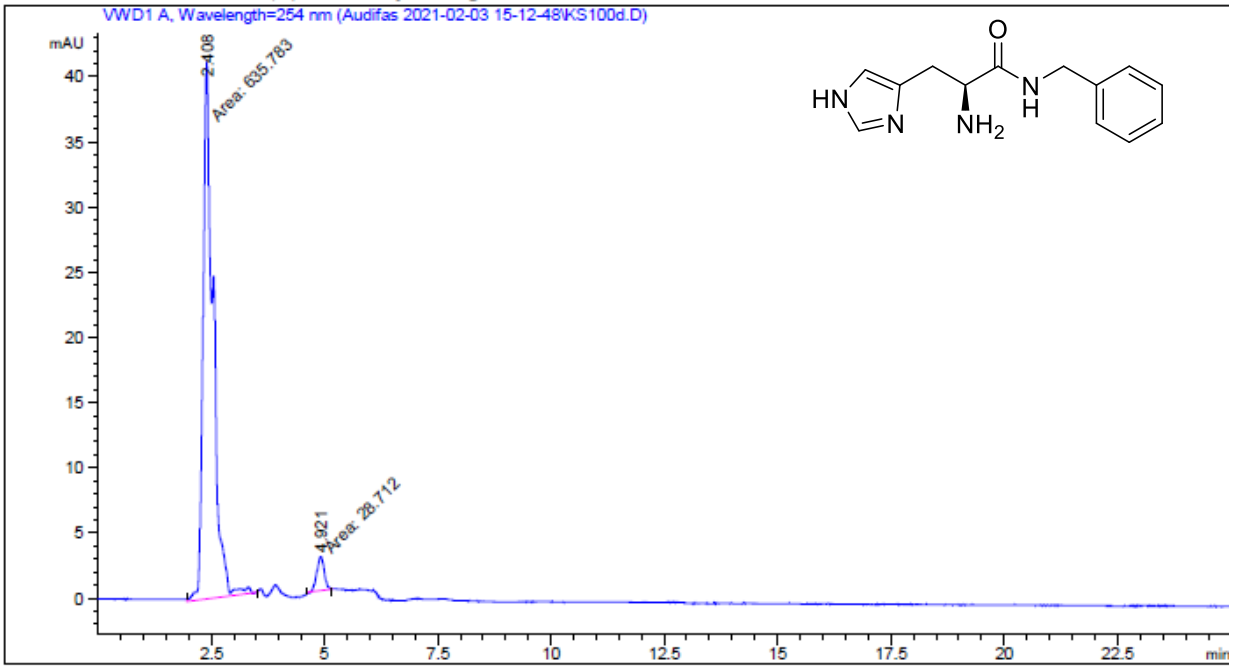
(*S*)-2-amino-*N*-benzyl-3-(1*H*-imidazol-4-yl)propanamide **(5a)**

```
Sample Name: KS100d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   3
Acq. Instrument : 1220 HPLC                       Location :   52
Injection Date  : 2/3/2021 4:09:20 PM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-02-03 15-12-48\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 2/3/2021 3:12:52 PM by SYSTEM
Method Info     : Methanol/Water 0.05 % TFA

Sample Info     : KS100d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.408 MM T  0.2569  635.78271   41.24117  95.6791
   2   4.921 MM T  0.2197   28.71196    2.57889   4.3209

Totals :                   664.49468   43.82006
```

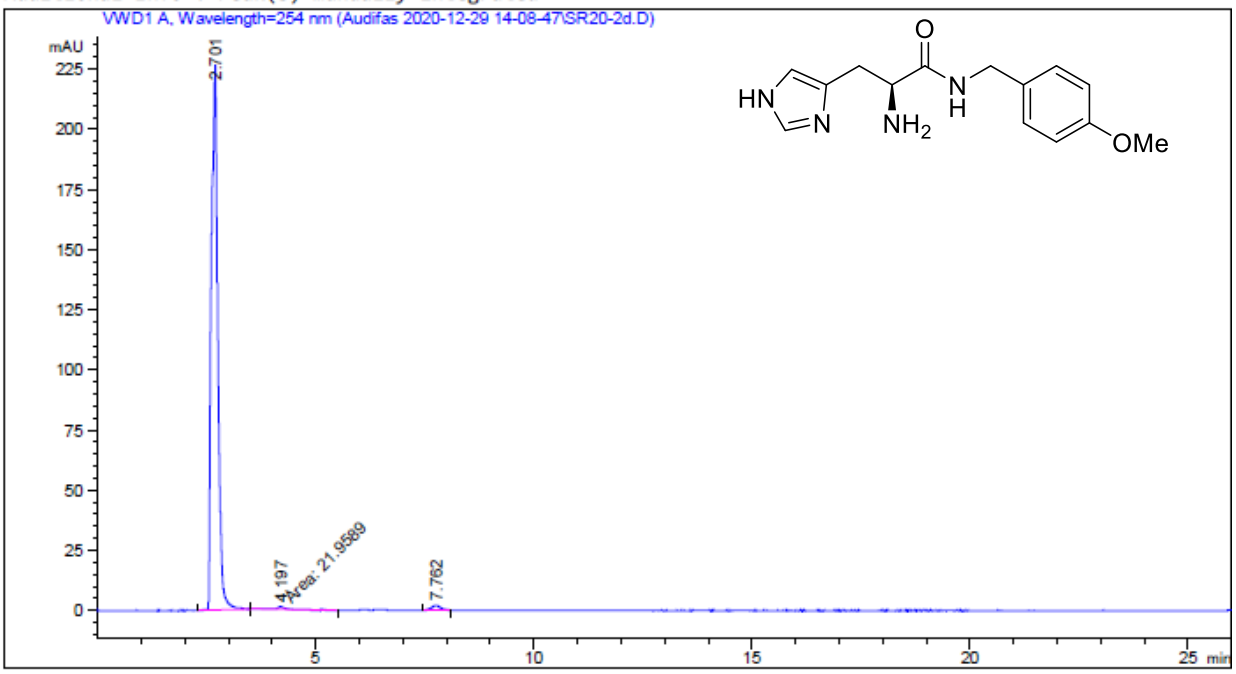
(*S*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(4-methoxybenzyl)propanamide **(5b)**

```
Sample Name: SR 20-2d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   7
Acq. Instrument : 1220 HPLC                       Location :   47
Injection Date  : 12/29/2020 5:01:24 PM              Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-29 14-08-47\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/29/2020 2:08:51 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05 %TFA / Methanol

Sample Info     : SR20-2d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs

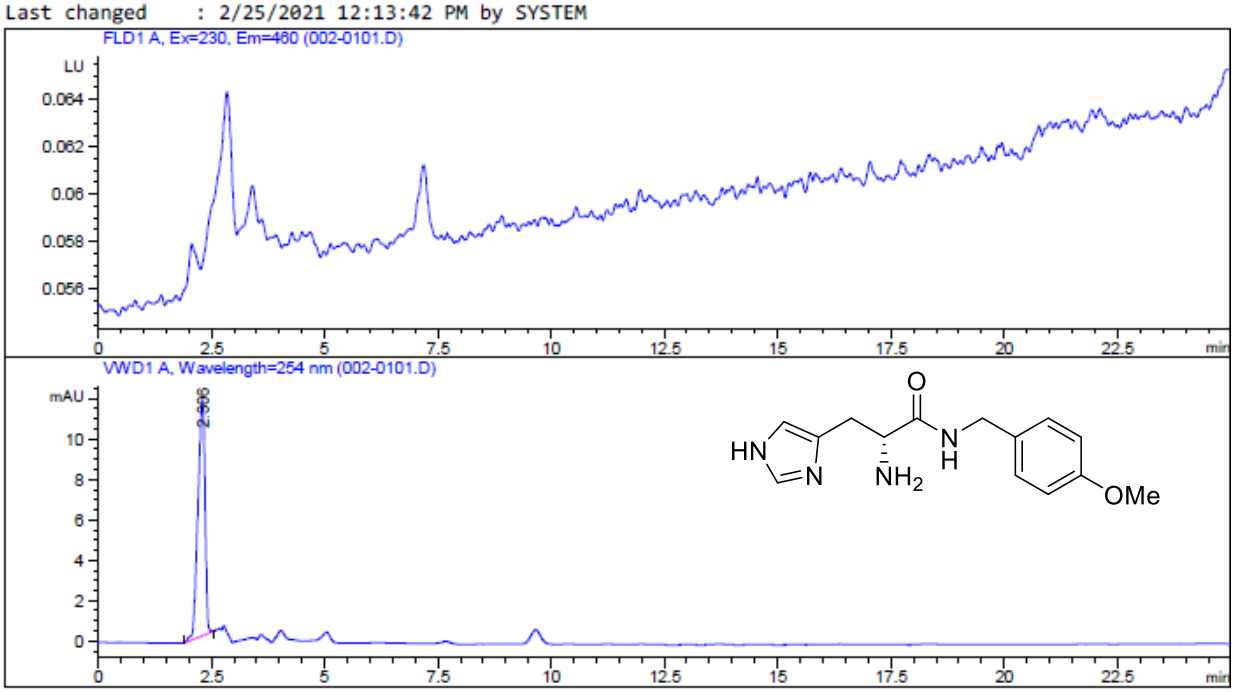

Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.701 BB    0.1384 2328.67993  226.55333  97.8858
   2   4.197 MM    0.3148   21.95889    1.16269   0.9230
   3   7.762 BB    0.1733   28.33842    1.92964   1.1912
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-methoxybenzyl)propanamide **(5c)**

```
Sample Name: PSK 12

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   1
Acq. Instrument : HPLC                           Location : Vial 2
Injection Date  : 2/25/2021 12:14:50 PM               Inj :   1
                                                 Inj Volume : 20.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2021-02-25 12-13-42\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2021-02-25 12-13-42\CQ.M (Sequence Method)
Last changed    : 2/25/2021 12:13:42 PM by SYSTEM
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=230, Em=460


Signal 2: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.306 BB    0.1743  127.70169   11.70746 100.0000

Totals :                   127.70169   11.70746
```

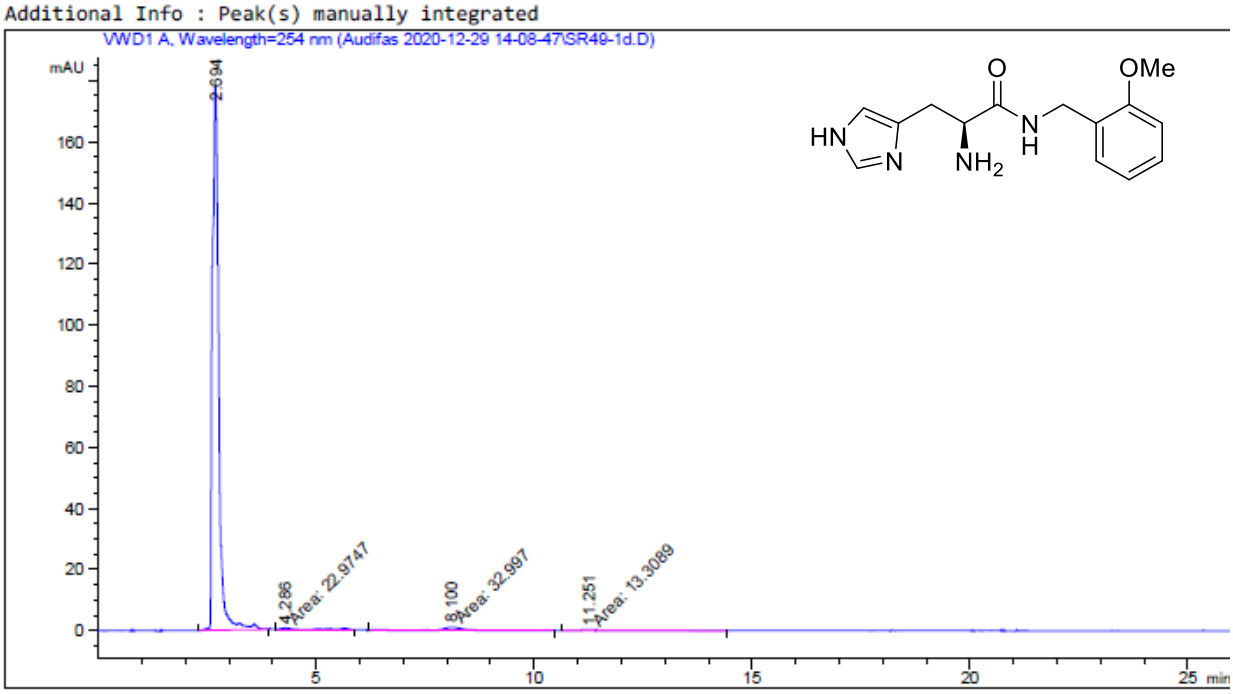
(*S*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(2-methoxybenzyl)propanamide **(5d)**

Sample Name: SR 48-1d

```
==================================================================
Acq. Operator   : SYSTEM                        Seq. Line :   2
Acq. Instrument : 1220 HPLC                      Location :   43
Injection Date  : 12/29/2020 2:38:59 PM              Inj :   1
                                                Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-29 14-08-47\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/29/2020 2:08:51 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05 %TFA / Methanol

Sample Info          SR48-1d

Additional Info : Peak(s) manually integrated
```




```
==================================================================
                      Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.694 BV R  0.1423 1776.13513  178.21967  96.2458
   2   4.286 MM    0.6600   22.97468 5.80142e-1   1.2450
   3   8.100 MM    0.5186   32.99696    1.06054   1.7881
   4  11.251 MM    1.2461   13.30886 1.78008e-1   0.7212
```

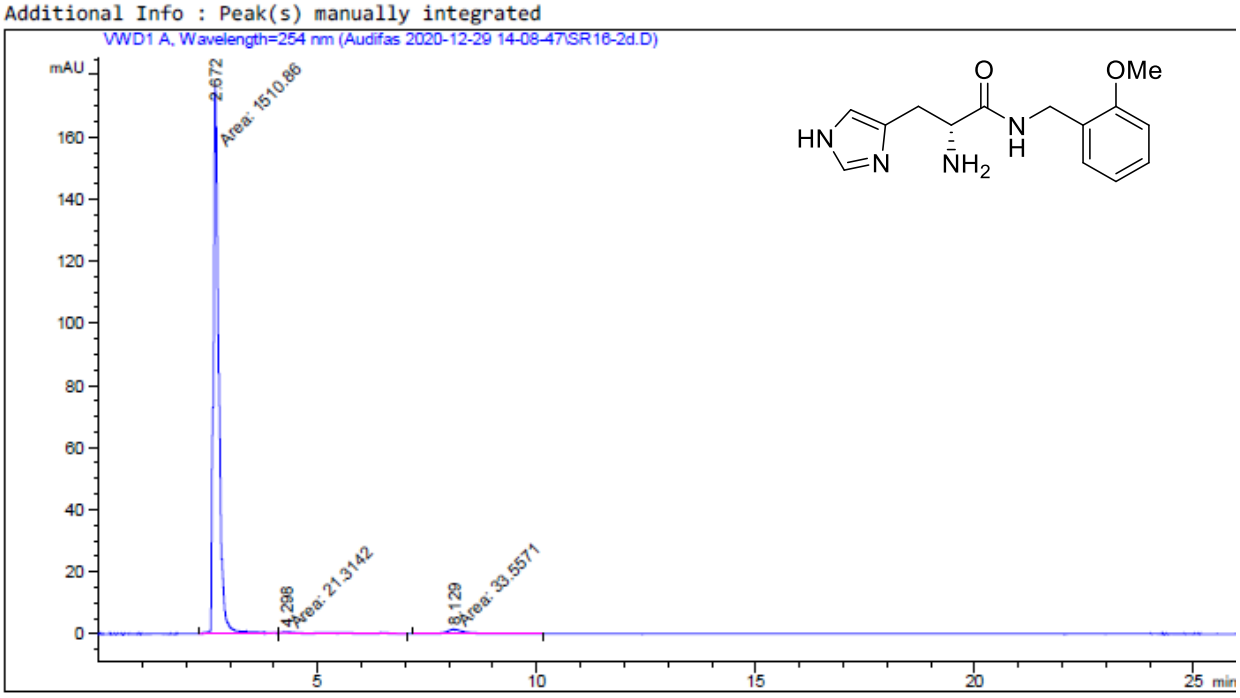
(*R*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(2-methoxybenzyl)propanamide **(5e)**

Sample Name: SR 16-2d

```
======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   8
Acq. Instrument : 1220 HPLC                        Location :   48
Injection Date  : 12/29/2020 5:29:52 PM                Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-29 14-08-47\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/29/2020 2:08:51 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05 %TFA / Methanol

Sample Info     : SR16-2d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.672 MM    0.1427 1510.85803  176.46837  96.4955
   2   4.298 MM    0.7663   21.31420 4.63570e-1   1.3613
   3   8.129 MM    0.4371   33.55715    1.27962   2.1432
```

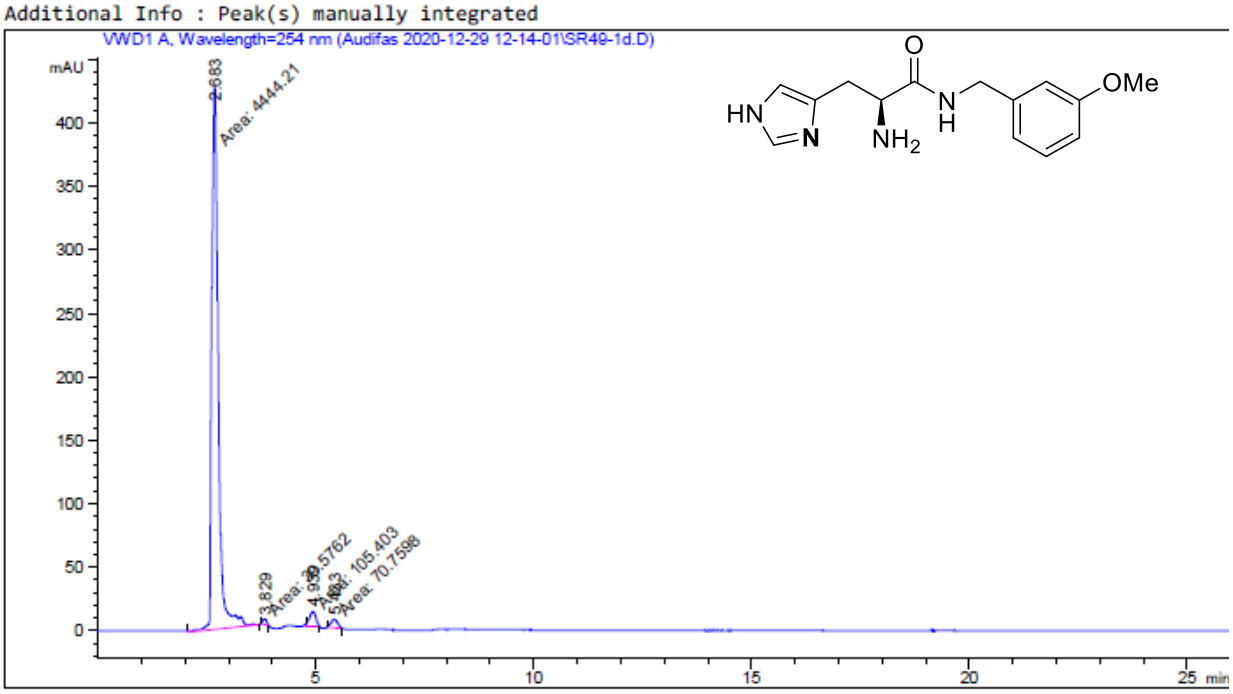
(*S*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(3-methoxybenzyl)propanamide **(5f)**

```
Sample Name: SR49-1d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   2
Acq. Instrument : 1220 HPLC                       Location :   41
Injection Date  : 12/29/2020 12:44:23 PM              Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-29 12-14-01\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/29/2020 12:14:04 PM by SYSTEM
Method Info     : Wash_Shafikur

Sample Info     : SR49-1d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs

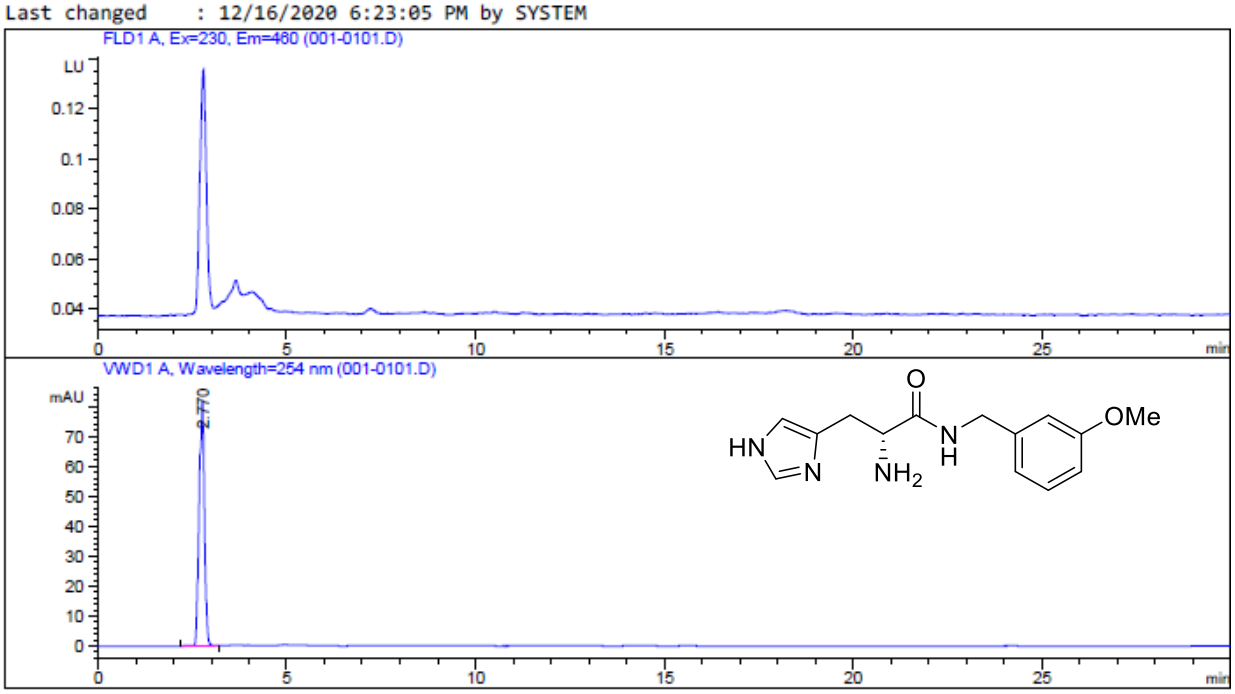

Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.683 MM T  0.3250 4444.20898  428.40399  95.5549
   2   3.829 MM    0.1060   30.57618    4.80739   0.6574
   3   4.939 MM    0.1619  105.40326   10.84810   2.2663
   4   5.433 MM    0.1813   70.75981    6.50518   1.5214
```

(*R*)-2-amino-3-(1*H*-3l4-imidazol-4-yl)-*N*-(3-methoxybenzyl)propanamide **(5g)**

```
Sample Name: SR 70-1 d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   1
Acq. Instrument : HPLC                           Location : Vial 1
Injection Date  : 12/16/2020 6:24:12 PM               Inj :   1
                                                 Inj Volume : 20.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2020-12-16 18-23-04\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2020-12-16 18-23-04\CQ.M (Sequence Method)
Last changed    : 12/16/2020 6:23:05 PM by SYSTEM
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=230, Em=460


Signal 2: VWD1 A, Wavelength=254 nm

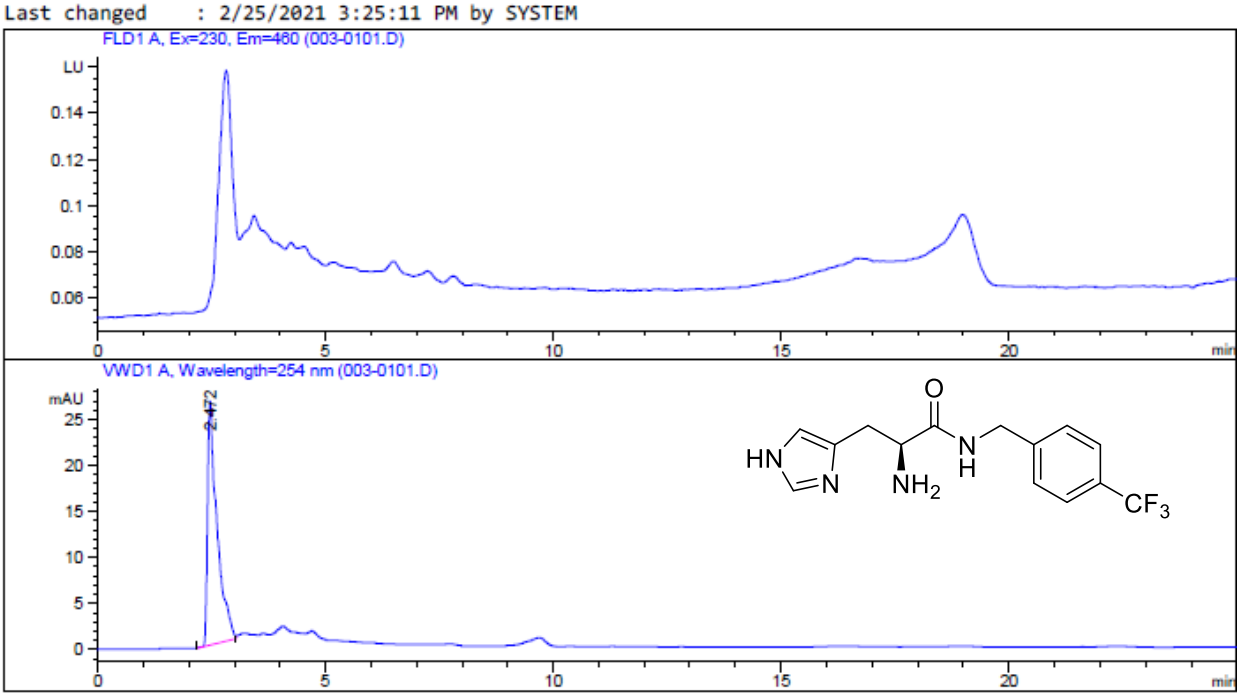
Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.770 BB    0.1274  790.94592   82.10738 100.0000

Totals :                   790.94592   82.10738
```

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)benzyl)propanamide **(5h)**

Sample Name: PSK 6d

```
======================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   1
Acq. Instrument : HPLC                         Location : Vial 3
Injection Date  : 2/25/2021 3:26:35 PM             Inj :   1
                                              Inj Volume : 50.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2021-02-25 15-25-11\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2021-02-25 15-25-11\CQ.M (Sequence Method)
Last changed    : 2/25/2021 3:25:11 PM by SYSTEM
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=230, Em=460

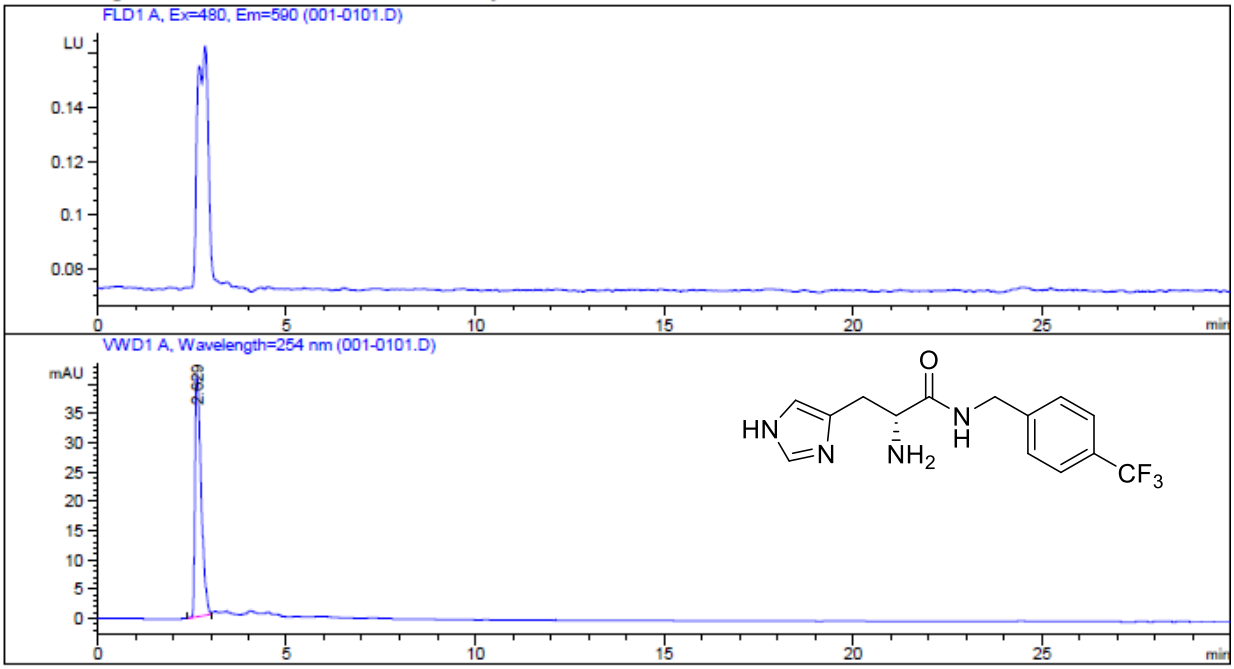

Signal 2: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.472 BV    0.1993  383.15048   26.40471 100.0000

Totals :                   383.15048   26.40471
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethyl)benzyl)propanamide **(5i)**

```
Sample Name: KS-102
========================================================================
Acq. Operator   : SYSTEM                        Seq. Line :   1
Acq. Instrument : HPLC                          Location : Vial 1
Injection Date  : 12/9/2020 12:08:59 PM              Inj :   1
                                               Inj Volume : 20.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2020-12-09 12-07-53\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2020-12-09 12-07-53\CQ.M (Sequence Method)
Last changed    : 12/9/2020 12:07:53 PM by SYSTEM
```




```
========================================================================
                        Area Percent Report
========================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=480, Em=590


Signal 2: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.629 BV    0.1720  464.86584   41.10150 100.0000

Totals :                   464.86584   41.10150
```

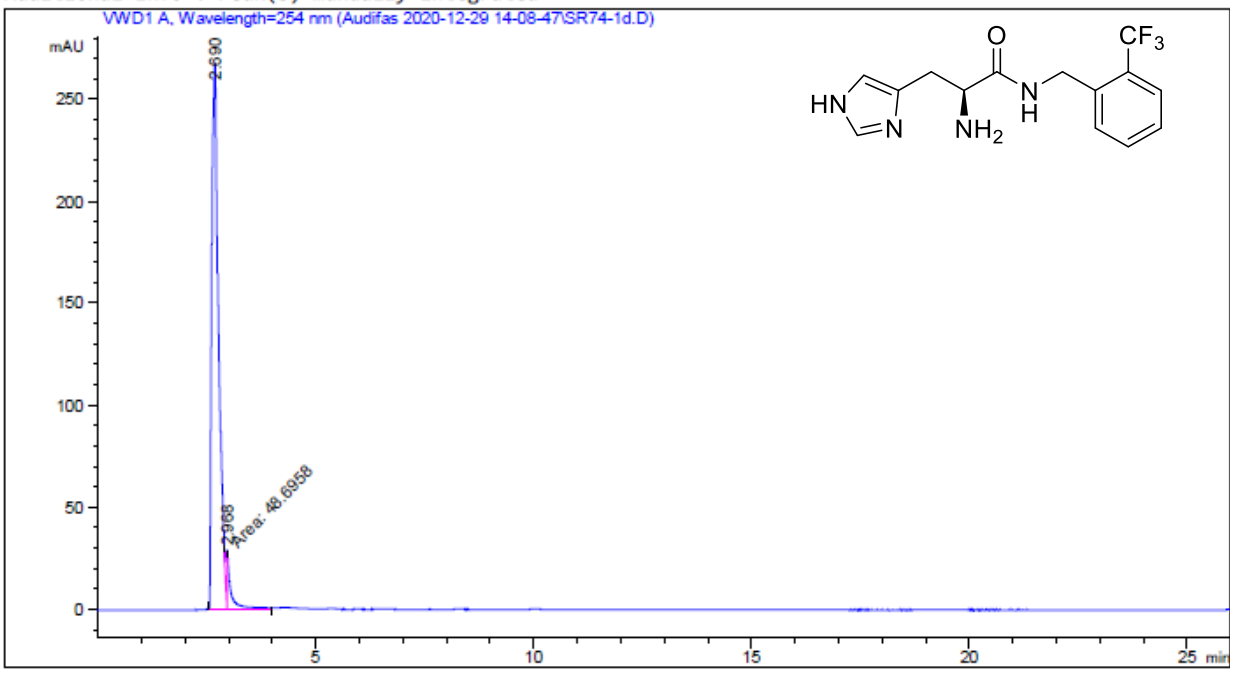
(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(trifluoromethyl)benzyl)propanamide **(5j)**

```
Sample Name: SR 74-1d

==================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   3
Acq. Instrument : 1220 HPLC                     Location :   44
Injection Date  : 12/29/2020 3:07:25 PM             Inj :   1
                                               Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-29 14-08-47\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/29/2020 2:08:51 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05 %TFA / Methanol

Sample Info     : SR74-1d

Additional Info : Peak(s) manually integrated
```




```
==================================================================
                      Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.690 MM R  0.1853 3100.65503  267.02853  98.4538
   2   2.968 MM T  0.0287   48.69577   28.29720   1.5462

Totals :                  3149.35080  295.32573
```

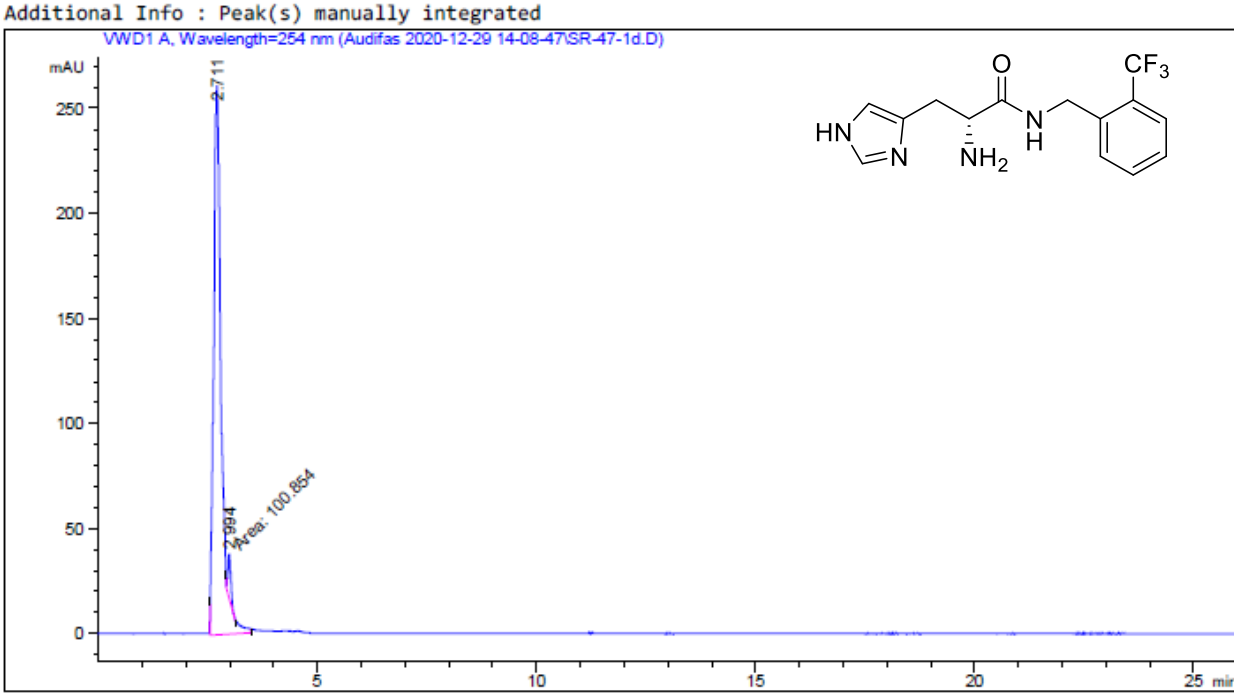
(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(trifluoromethyl)benzyl)propanamide

```
Sample Name: SR 47-1d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :  30
Acq. Instrument : 1220 HPLC                       Location :   64
Injection Date  : 12/30/2020 4:03:14 AM               Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-29 14-08-47\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/29/2020 2:08:51 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05 %TFA / Methanol

Sample Info     : SR-47-1d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.711 MM R  0.1988 3119.55469  261.53845  96.8683
   2   2.994 MM T  0.0867  100.85400   21.43935   3.1317

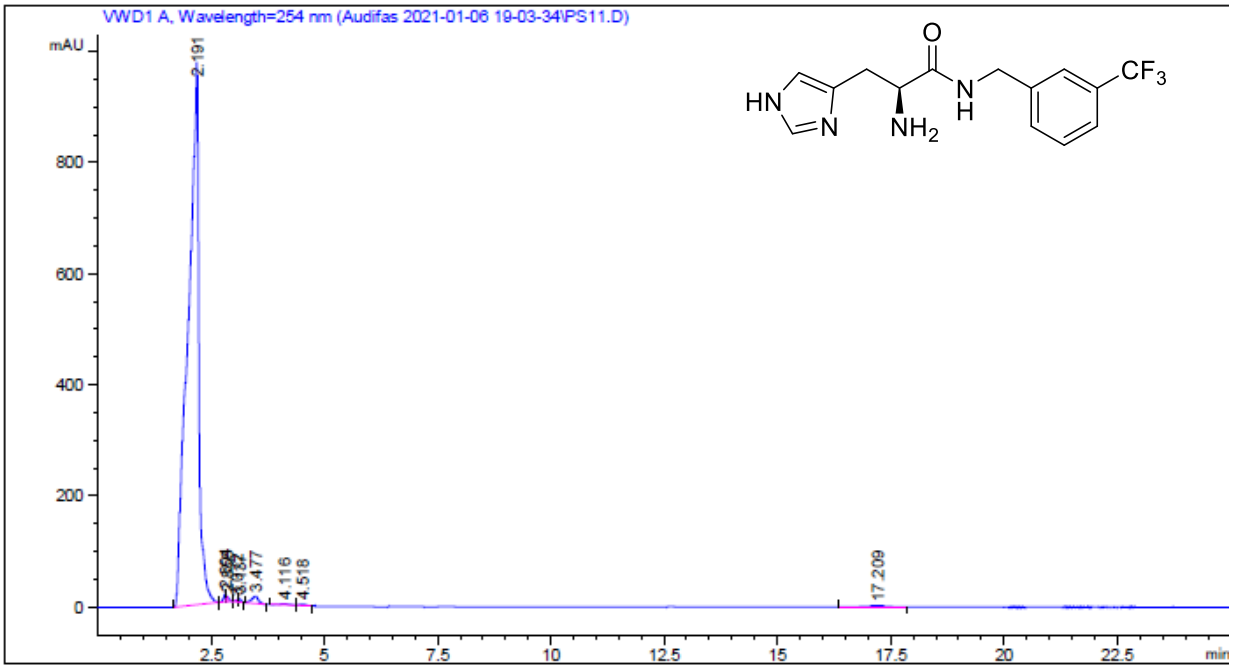
Totals :                  3220.40869  282.97780
```

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-(trifluoromethyl)benzyl)propanamide **(5l)**

```
Sample Name: PS11

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :  33
Acq. Instrument : 1220 HPLC                       Location :   83
Injection Date  : 1/7/2021 9:50:42 AM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : PS11
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]       [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.191 BV    0.2070 1.60680e4   976.16382  97.6645
   2   2.804 VV    0.0523   50.54400   13.62504   0.3072
   3   2.856 VB    0.0494   31.15095    9.23310   0.1893
   4   3.072 BV    0.0430    7.76413    2.20815   0.0472
   5   3.137 VB    0.0726   22.62145    4.70742   0.1375
```

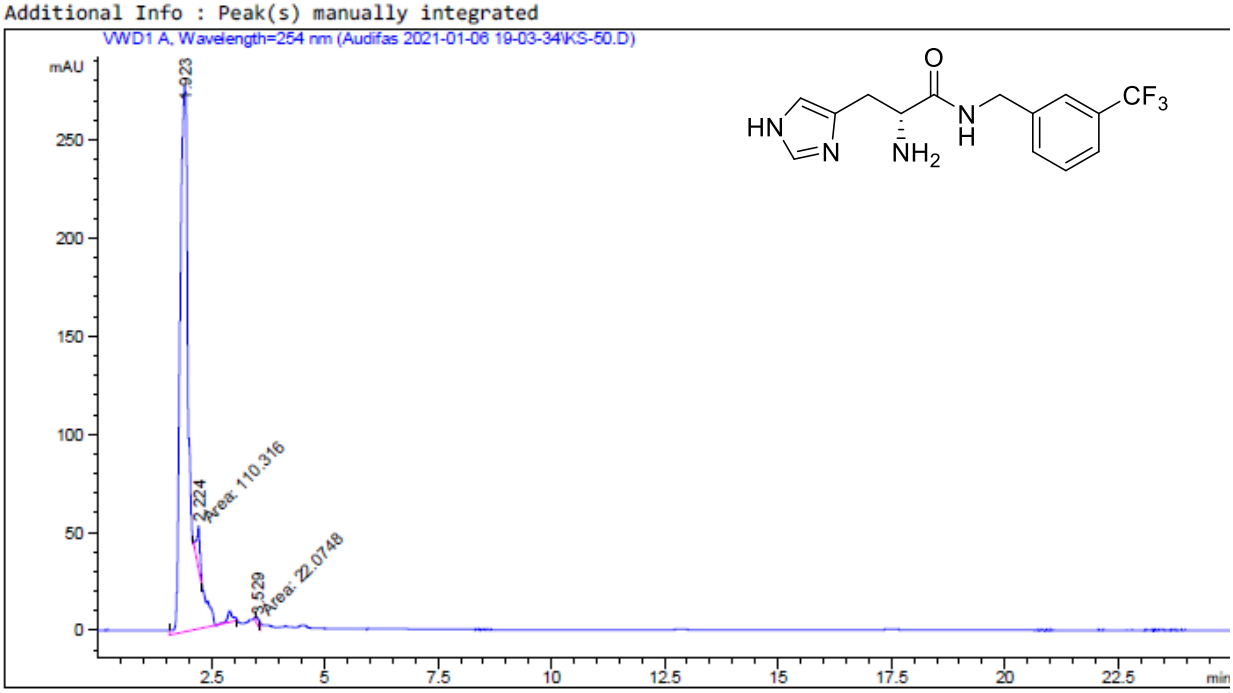
(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-(trifluoromethyl)benzyl)propanamide **(5m)**

```
Sample Name: KS-50

==================================================================
Acq. Operator   : SYSTEM                       Seq. Line :  24
Acq. Instrument : 1220 HPLC                    Location :   76
Injection Date  : 1/7/2021 5:42:46 AM               Inj :   1
                                              Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : KS-50

Additional Info : Peak(s) manually integrated
```




```
==================================================================
                      Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

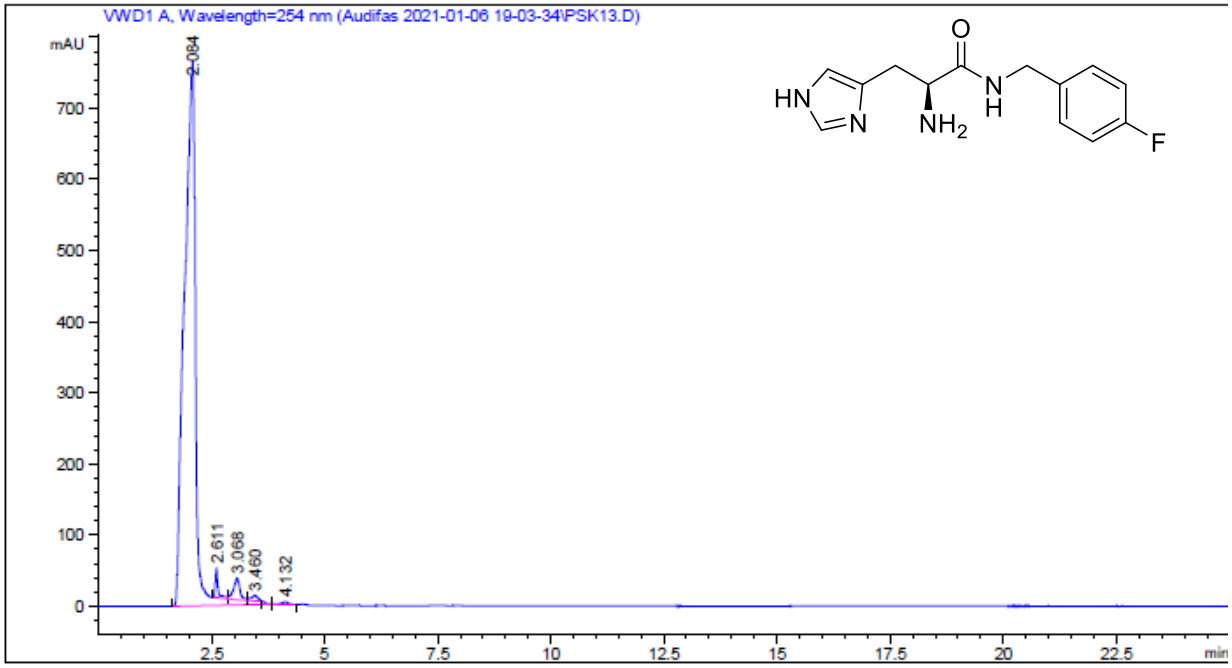
Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.923 MM R  0.2843 4027.58911  278.71985  96.8175
   2   2.224 MM T  0.1064  110.31607   22.17361   2.6518
   3   3.529 MM T  0.0739   22.07477    3.53974   0.5306
```

(*S*)-2-amino-*N*-(4-fluorobenzyl)-3-(1*H*-imidazol-4-yl)propanamide **(5n)**

```
Sample Name: PSK13

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :  25
Acq. Instrument : 1220 HPLC                       Location :   77
Injection Date  : 1/7/2021 6:10:19 AM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water  0.05 %

Sample Info     : PSK13
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.084 BV R  0.2349 1.32136e4   765.27478  95.7218
   2   2.611 VV E  0.0615  169.56122   40.40717   1.2283
   3   3.068 VV E  0.1376  300.43854   30.57756   2.1764
   4   3.460 VB E  0.1309   78.70617    7.52999   0.5702
   5   4.132 VV R  0.1410   41.86796    3.52590   0.3033
```

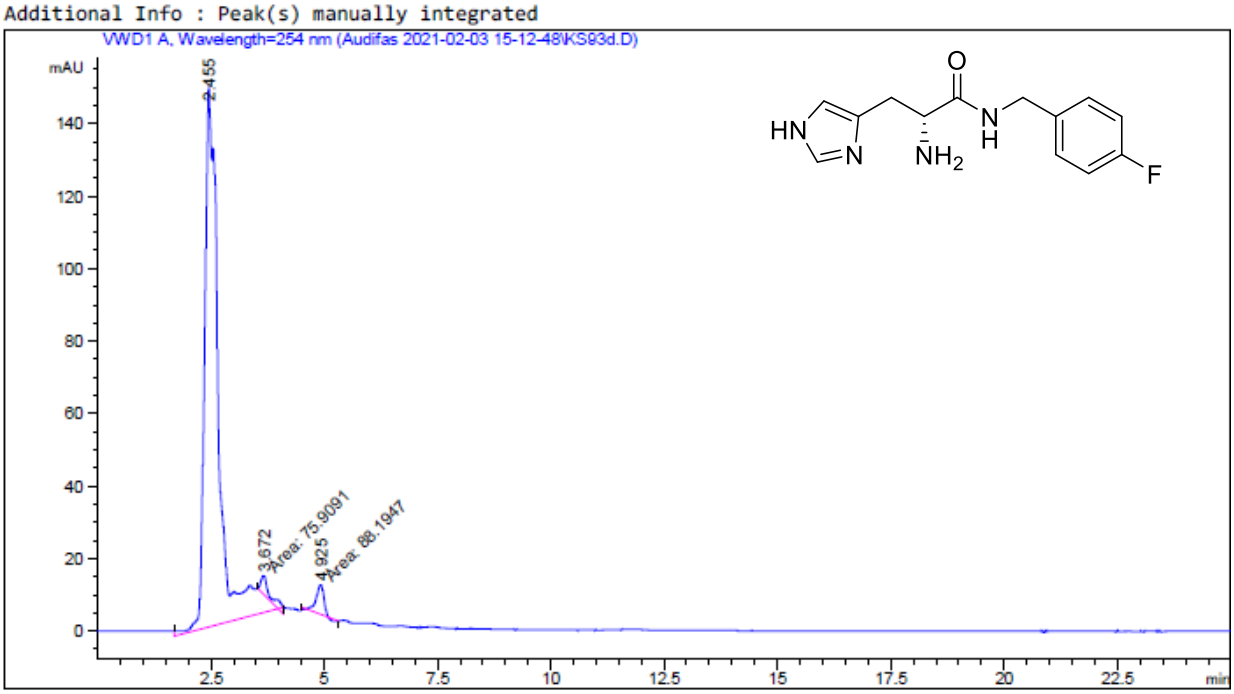
(*R*)-2-amino-*N*-(4-fluorobenzyl)-3-(1*H*-imidazol-4-yl)propanamide **(5o)**

```
Sample Name: KS 93d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   4
Acq. Instrument : 1220 HPLC                       Location :   53
Injection Date  : 2/3/2021 4:36:50 PM                  Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-02-03 15-12-48\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 2/3/2021 3:12:52 PM by SYSTEM
Method Info     : Methanol/Water 0.05 % TFA

Sample Info     : KS93d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs

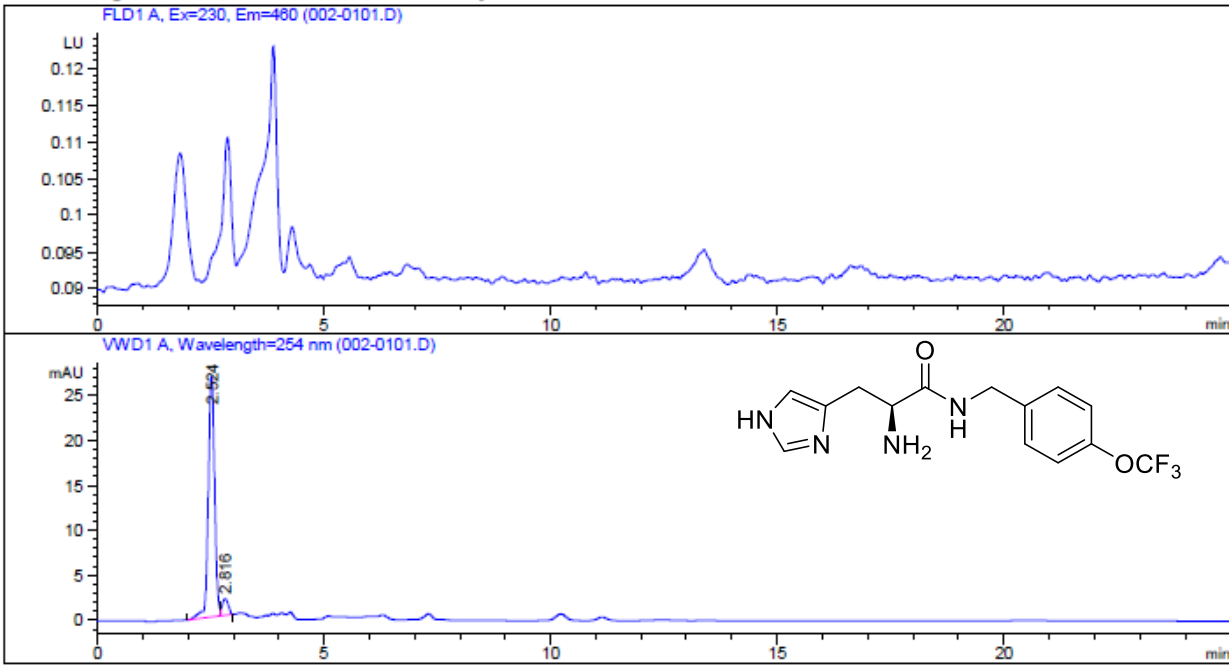

Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.455 MM R  0.3552 3183.49414  149.36600  95.0979
   2   3.672 MM T  0.2428   75.90913    5.21036   2.2676
   3   4.925 MM T  0.1811   88.19469    8.11711   2.6346
```

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethoxy)benzyl)propanamide **(5p)**

```
Sample Name: PSK 5d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   1
Acq. Instrument : HPLC                           Location : Vial 2
Injection Date  : 2/22/2021 1:15:35 PM                Inj :   1
                                                 Inj Volume : 20.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2021-02-22 13-14-27\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2021-02-22 13-14-27\CQ.M (Sequence Method)
Last changed    : 2/22/2021 1:14:27 PM by SYSTEM
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=230, Em=460


Signal 2: VWD1 A, Wavelength=254 nm

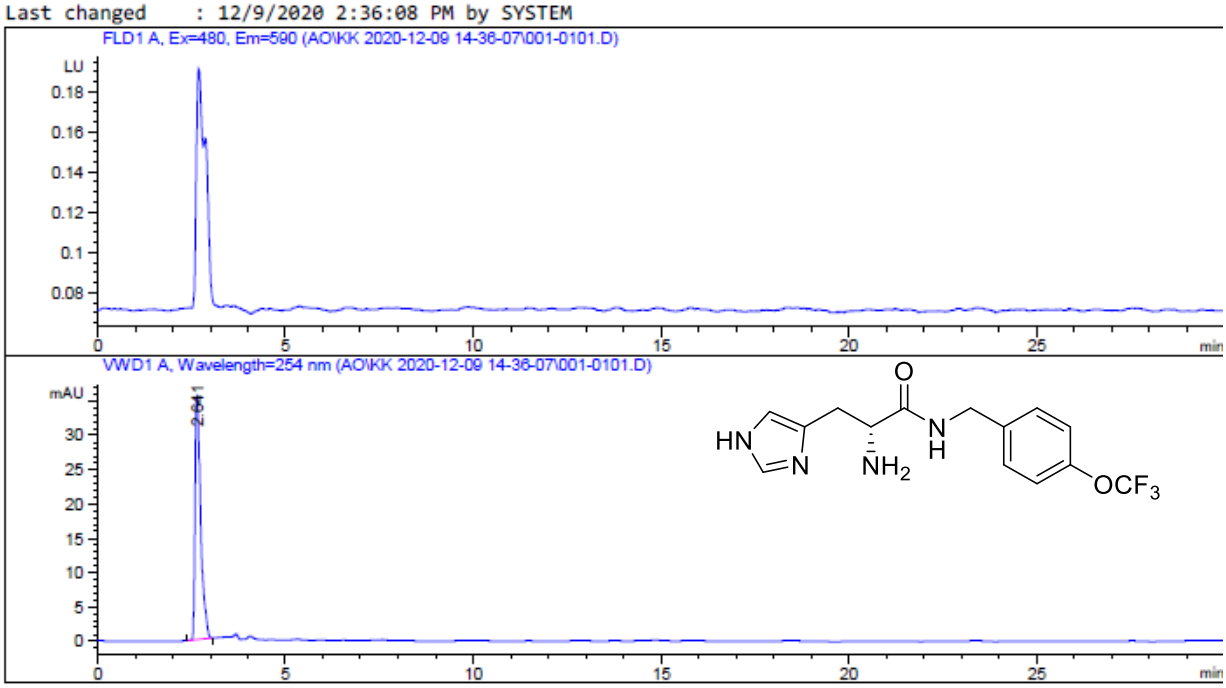
Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.524 BV    0.1479  261.86145   26.82778  93.8184
   2   2.816 VB    0.1470   17.25385    1.84755   6.1816

Totals :                   279.11530   28.67532
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-(trifluoromethoxy)benzyl)propanamide **(5q)**

Sample Name: KS-99

```
==================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   1
Acq. Instrument : HPLC                          Location : Vial 1
Injection Date  : 12/9/2020 2:37:16 PM              Inj :   1
                                               Inj Volume : 20.000 µl
Method          : C:\CHEM32\1\DATA\AO\KK 2020-12-09 14-36-07\CQ.M (Sequence Method)
Last changed    : 12/9/2020 2:36:08 PM by SYSTEM
```




```
==================================================================
                        Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=480, Em=590


Signal 2: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.641 BB    0.1623  368.46179   35.42866 100.0000

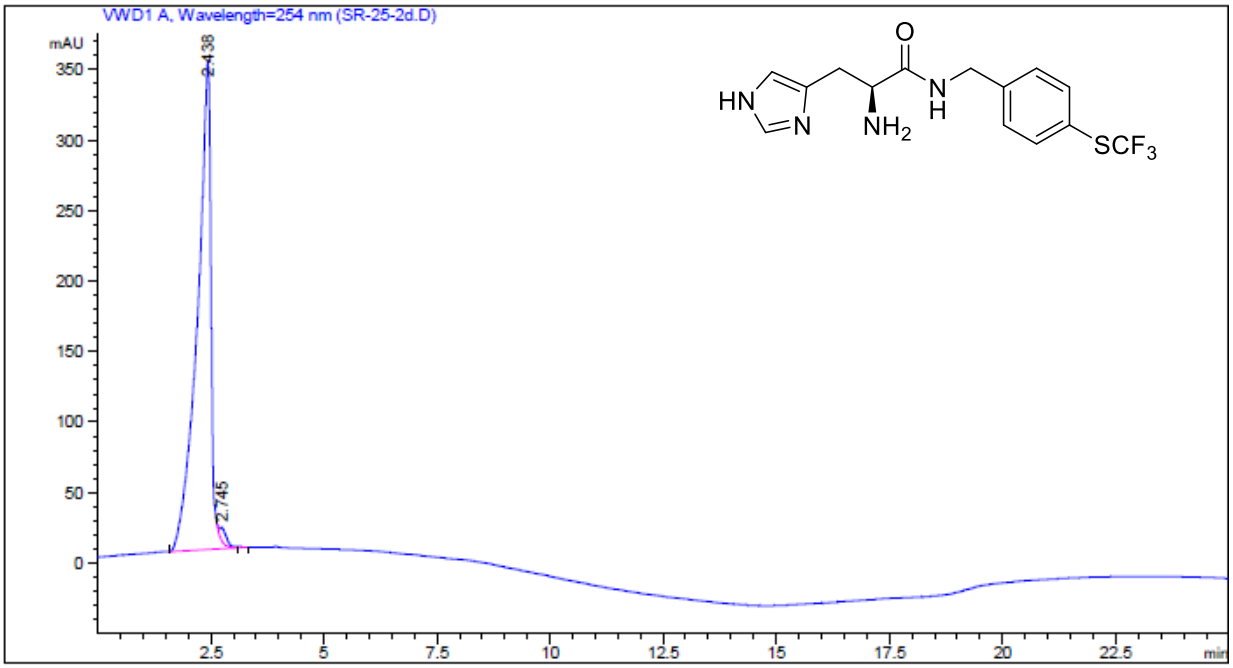
Totals :                   368.46179   35.42866
```

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-((trifluoromethyl)thio)benzyl)propanamide

Sample Name: SR-25-2d

```
==================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   6
Acq. Instrument : 1220 HPLC                     Location :   56
Injection Date  : 12/30/2020 5:50:10 PM             Inj :   1
                                               Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Sequence File   : C:\Chem32\1\Data\Audifas 2020-12-30 15-31-06\Audifas.S
Method          : C:\Chem32\1\Data\Audifas 2020-12-30 15-31-06\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/30/2020 3:31:09 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05% TFA / Methanol

Sample Info     : SR-25-2d
```




```
==================================================================
                      Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.438 BV R  0.2771 7274.41016  346.24774  98.5591
   2   2.745 VV E  0.1413  106.35146    9.52277   1.4409

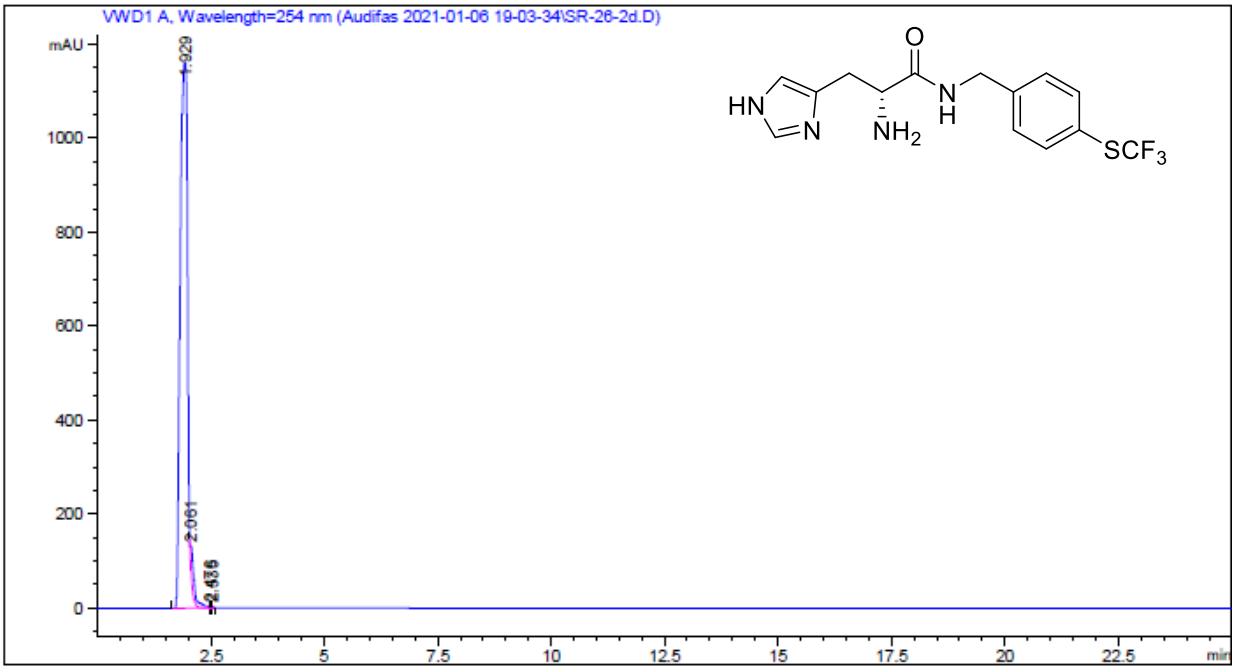
Totals :                  7380.76161  355.77051
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-((trifluoromethyl)thio)benzyl)propanamide

Sample Name: SR-26-2d

```
==================================================================
Acq. Operator   : SYSTEM                       Seq. Line :  11
Acq. Instrument : 1220 HPLC                     Location :  67
Injection Date  : 1/6/2021 11:44:37 PM              Inj :  1
                                               Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR-26-2d
```




```
==================================================================
                        Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs

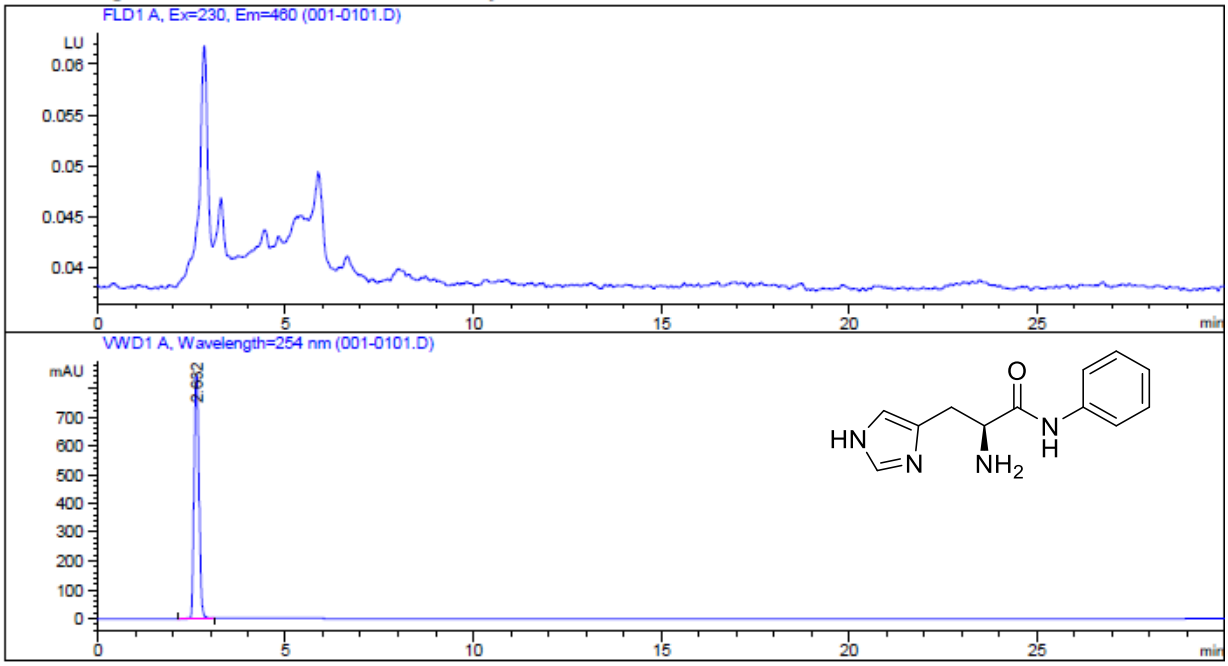

Signal 1: VWD1 A, Wavelength=254 nm
```

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	1.929	BV R	0.1848	1.33215e4	1161.39929	96.5065
2	2.061	VV E	0.1207	467.32349	50.17341	3.3855
3	2.476	VV E	0.0413	8.78115	2.94324	0.0636
4	2.535	VB E	0.0416	6.12294	2.15543	0.0444

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenylpropanamide **(6a)**

```
Sample Name: SR 73-1 d

======================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   1
Acq. Instrument : HPLC                         Location : Vial 1
Injection Date  : 12/16/2020 4:40:42 PM            Inj :   1
                                              Inj Volume : 20.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2020-12-16 16-39-32\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2020-12-16 16-39-32\CQ.M (Sequence Method)
Last changed    : 12/16/2020 4:39:33 PM by SYSTEM
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=230, Em=460


Signal 2: VWD1 A, Wavelength=254 nm

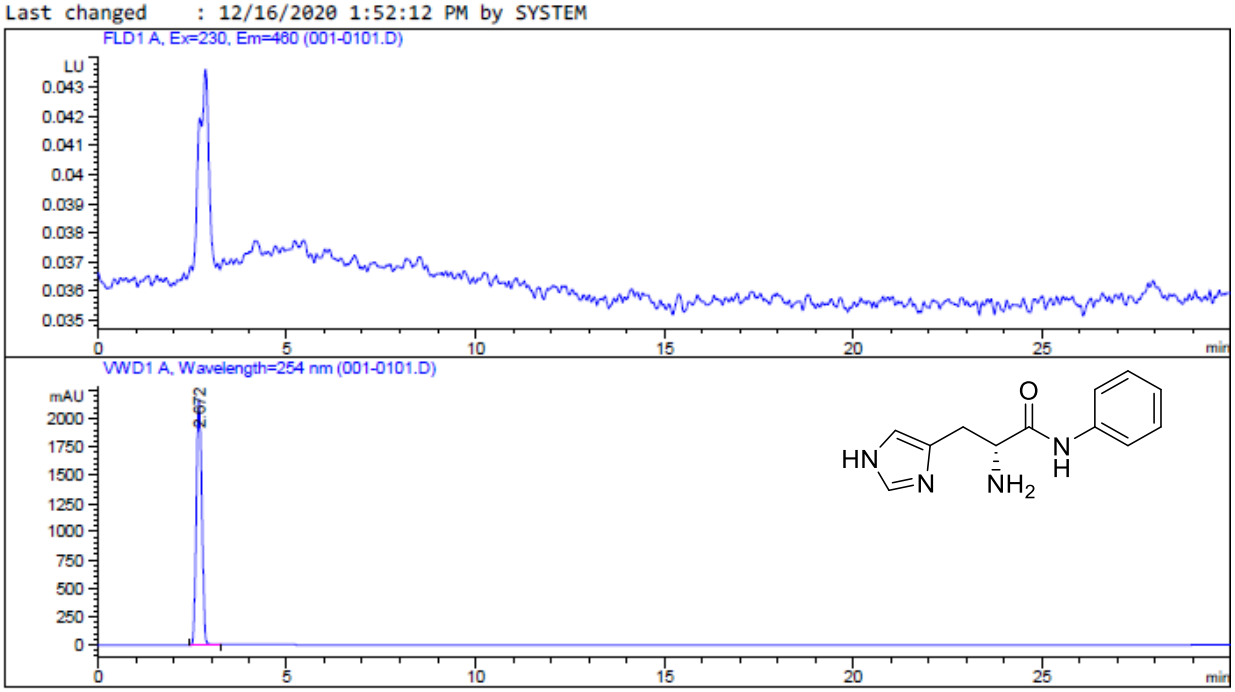
Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.632 BB    0.1416 7673.69678  848.12286 100.0000

Totals :                  7673.69678  848.12286
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-phenylpropanamide **(6b)**

```
Sample Name: SR 72-1 d

======================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   1
Acq. Instrument : HPLC                         Location : Vial 1
Injection Date  : 12/16/2020 1:53:19 PM             Inj :   1
                                              Inj Volume : 20.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2020-12-16 13-52-11\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2020-12-16 13-52-11\CQ.M (Sequence Method)
Last changed    : 12/16/2020 1:52:12 PM by SYSTEM
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs


Signal 1: FLD1 A, Ex=230, Em=460


Signal 2: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.672 BB    0.1621 2.19906e4  2173.03809 100.0000

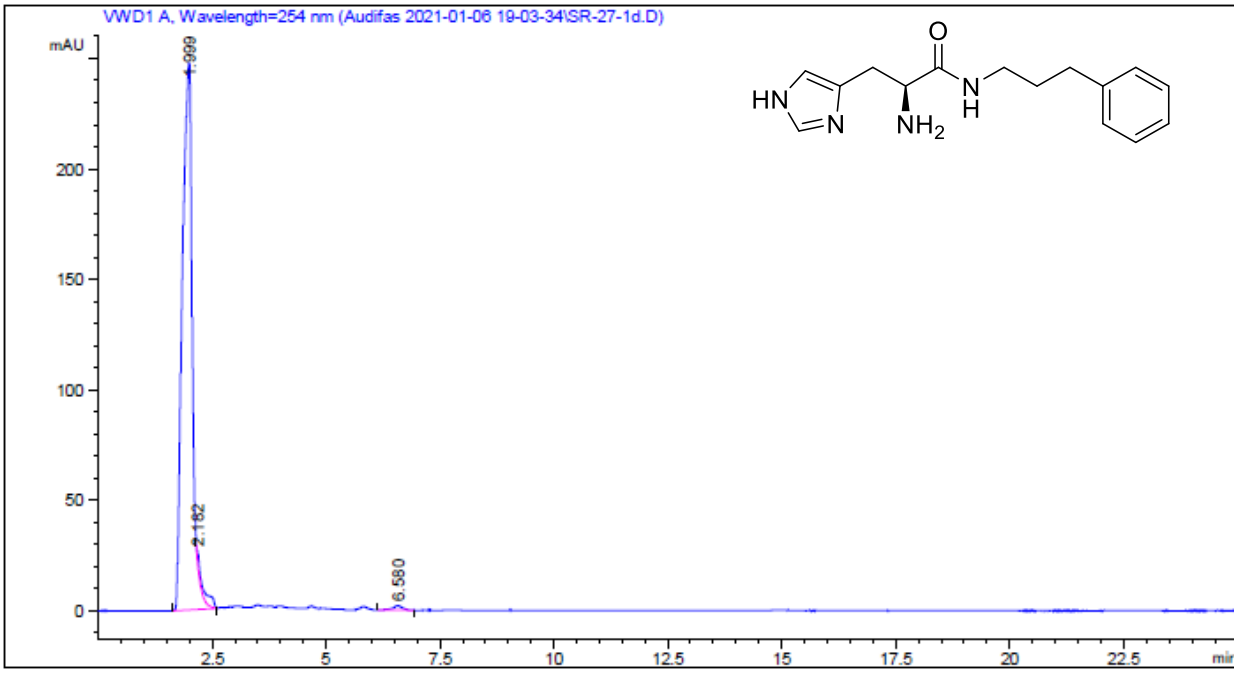
Totals :                  2.19906e4  2173.03809
```

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-phenylpropyl)propanamide **(7a)**

```
Sample Name: SR-27-1d

=====================================================================
Acq. Operator   : SYSTEM                          Seq. Line :  18
Acq. Instrument : 1220 HPLC                        Location :  72
Injection Date  : 1/7/2021 2:57:24 AM                  Inj :  1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR-27-1d
```




```
=====================================================================
                        Area Percent Report
=====================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.999 BV R  0.2161 3774.33862  247.89630  96.7906
   2   2.182 VB E  0.2153   96.46211    5.44194   2.4737
   3   6.580 VB    0.1799   28.68609    1.88636   0.7356

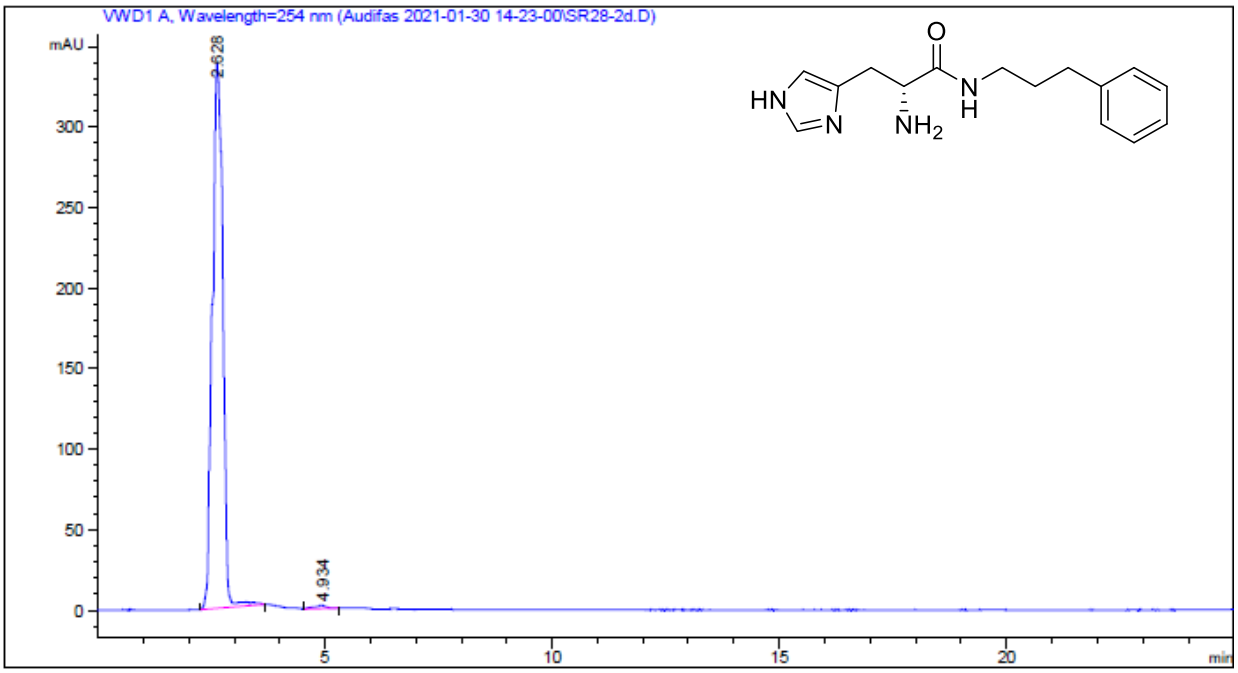
Totals :                  3899.48682  255.22460
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(3-phenylpropyl)propanamide **(7b)**

```
Sample Name: SR28-2d

======================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   3
Acq. Instrument : 1220 HPLC                    Location :   53
Injection Date  : 1/30/2021 3:19:52 PM             Inj :   1
                                              Inj Volume : 50.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-30 14-23-00\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/30/2021 2:23:03 PM by SYSTEM
Method Info     : Methanol/Water 0.05 % TFA

Sample Info     : SR28-2d
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.628 BV R  0.2147 5423.52588  338.86234  99.3150
   2   4.934 BB    0.2309   37.40524    1.90636   0.6850

Totals :                  5460.93111  340.76869
```

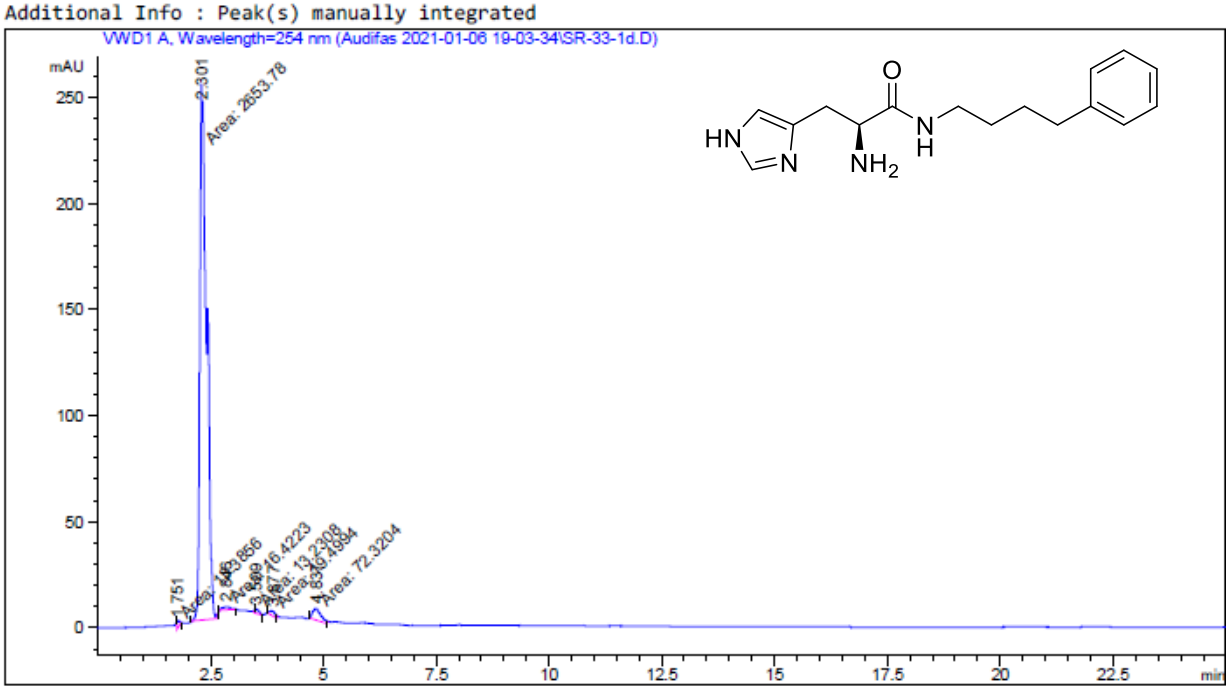
(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-phenylbutyl)propanamide **(8a)**

```
Sample Name: SR-33-1d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :  20
Acq. Instrument : 1220 HPLC                        Location :  74
Injection Date  : 1/7/2021 3:52:26 AM                  Inj :  1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR-33-1d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm
```

Peak #	RetTime [min]	Type	Width [min]	Area [mAU*s]	Height [mAU]	Area %
1	1.751	MM T	0.0477	13.38564	3.50553	0.4800
2	2.301	MM T	0.1753	2653.78223	252.33617	95.1640
3	2.846	MM T	0.1871	16.42233	1.46263	0.5889
4	3.509	MM T	0.1108	13.23078	1.99102	0.4745

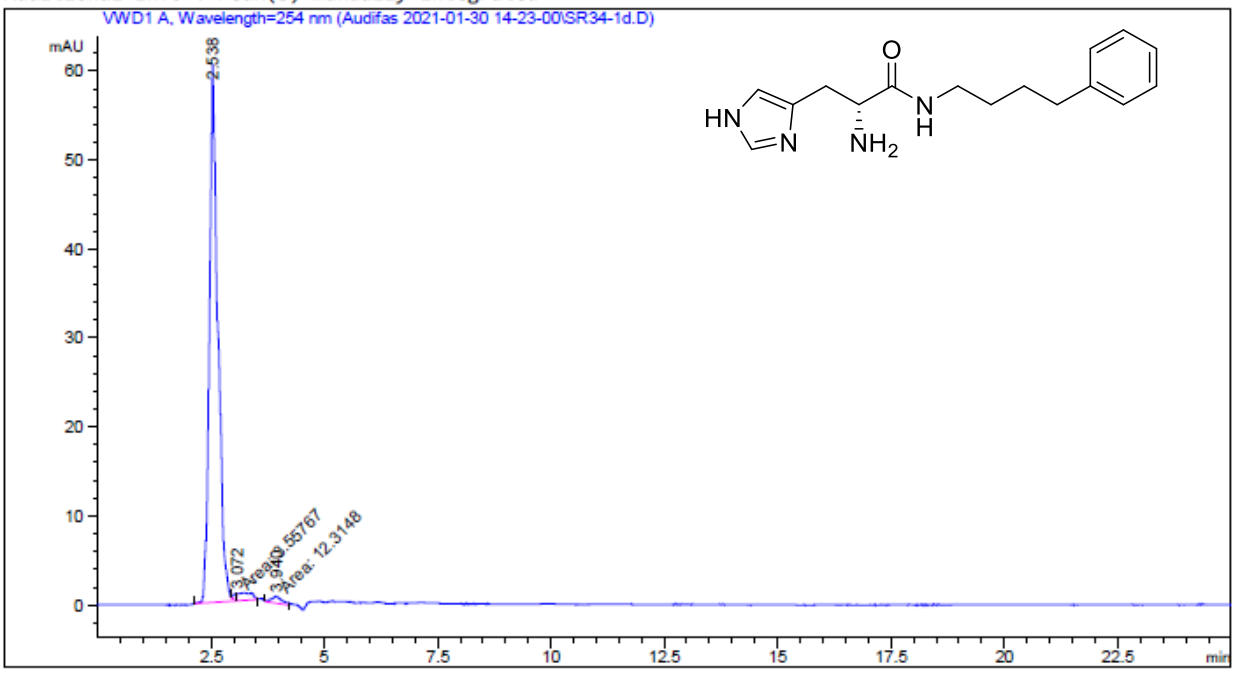
(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(4-phenylbutyl)propanamide **(8b)**

```
Sample Name: SR34-1d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   1
Acq. Instrument : 1220 HPLC                       Location :   51
Injection Date  : 1/30/2021 2:24:48 PM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-30 14-23-00\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/30/2021 2:23:03 PM by SYSTEM
Method Info     : Methanol/Water 0.05 % TFA

Sample Info     : SR34-1d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.538 MM R  0.2788  802.77618   60.50960  98.0611
   2   3.072 MM T  0.0814    3.55767 7.28136e-1   0.4346
   3   3.940 MM T  0.2576   12.31482 7.74216e-1   1.5043
```

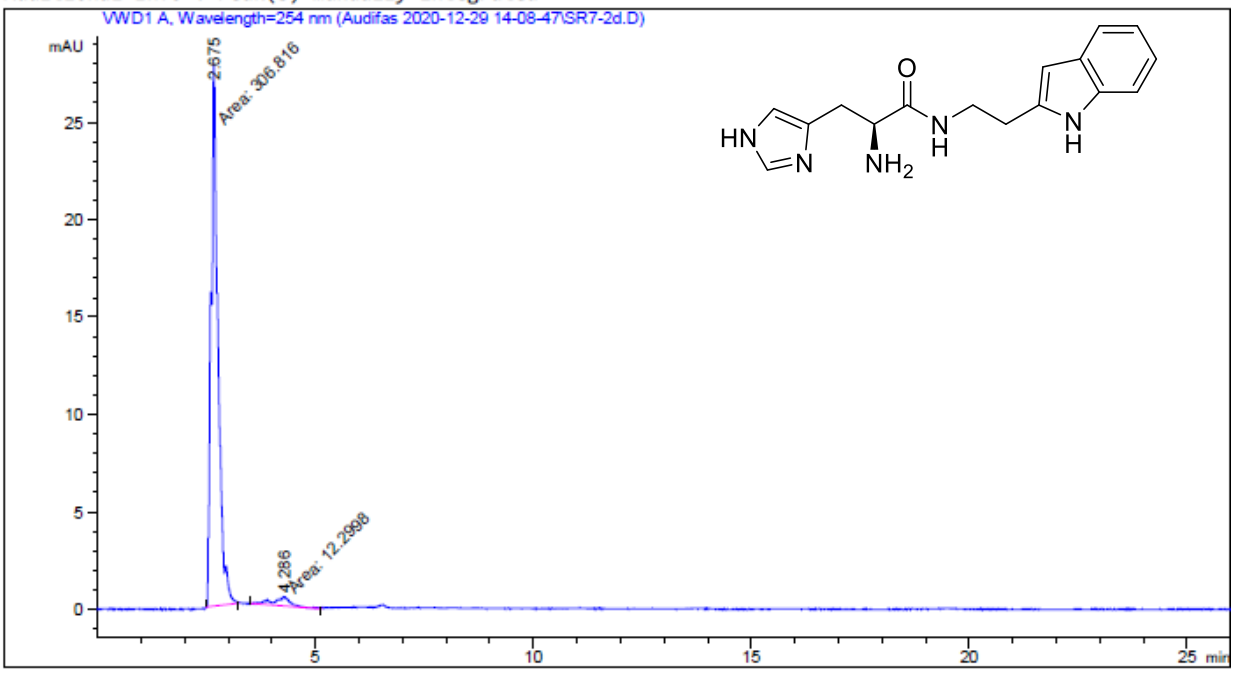
(*S*)-*N*-(2-(1*H*-indol-3-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide **(9d)**

```
Sample Name: SR7-2d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   4
Acq. Instrument : 1220 HPLC                       Location :   45
Injection Date  : 12/29/2020 3:35:52 PM               Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-29 14-08-47\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/29/2020 2:08:51 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05 %TFA / Methanol

Sample Info     : SR7-2d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.675 MM    0.1840  306.81577   27.78871  96.1457
   2   4.286 MM    0.4499   12.29981 4.55628e-1   3.8543

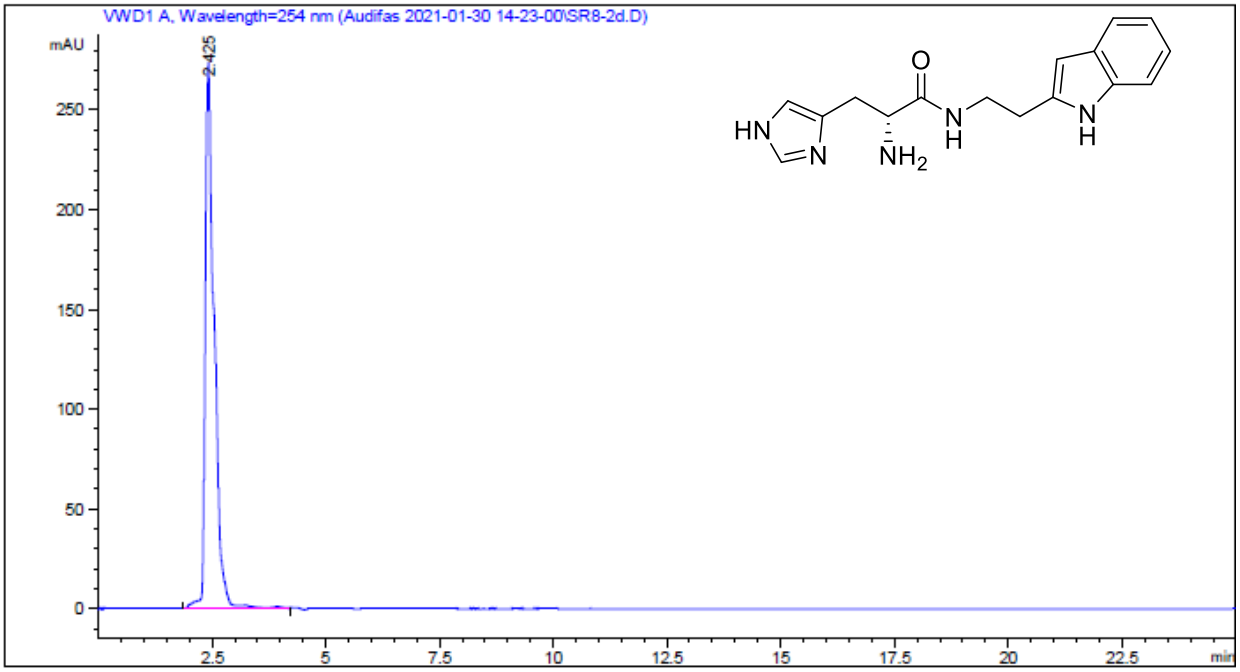
Totals :                   319.11558   28.24434
```

(*R*)-*N*-(2-(1*H*-indol-3-yl)ethyl)-2-amino-3-(1*H*-imidazol-4-yl)propanamide **(9e)**

```
Sample Name: SR8-2d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   4
Acq. Instrument : 1220 HPLC                       Location :   54
Injection Date  : 1/30/2021 3:47:27 PM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-30 14-23-00\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/30/2021 2:23:03 PM by SYSTEM
Method Info     : Methanol/Water 0.05 % TFA

Sample Info     : SR8-2d
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.425 BV R  0.1908 3754.30176  274.04691 100.0000

Totals :                  3754.30176  274.04691
```

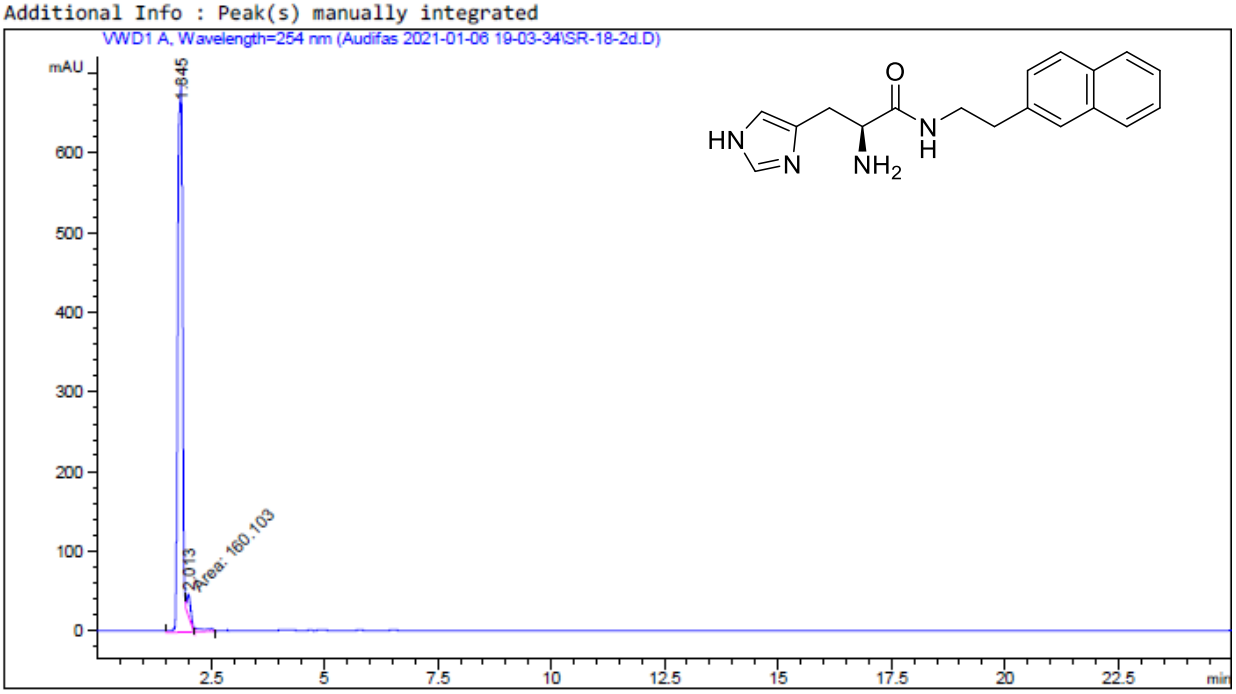
(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(naphthalen-2-yl)ethyl)propanamide **(9f)**

```
Sample Name: SR-18-2d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   6
Acq. Instrument : 1220 HPLC                       Location :   64
Injection Date  : 1/6/2021 9:26:47 PM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR-18-2d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.845 MM R  0.1948 5318.42285  687.43878  97.0776
   2   2.013 MM T  0.1050  160.10272   26.58394   2.9224

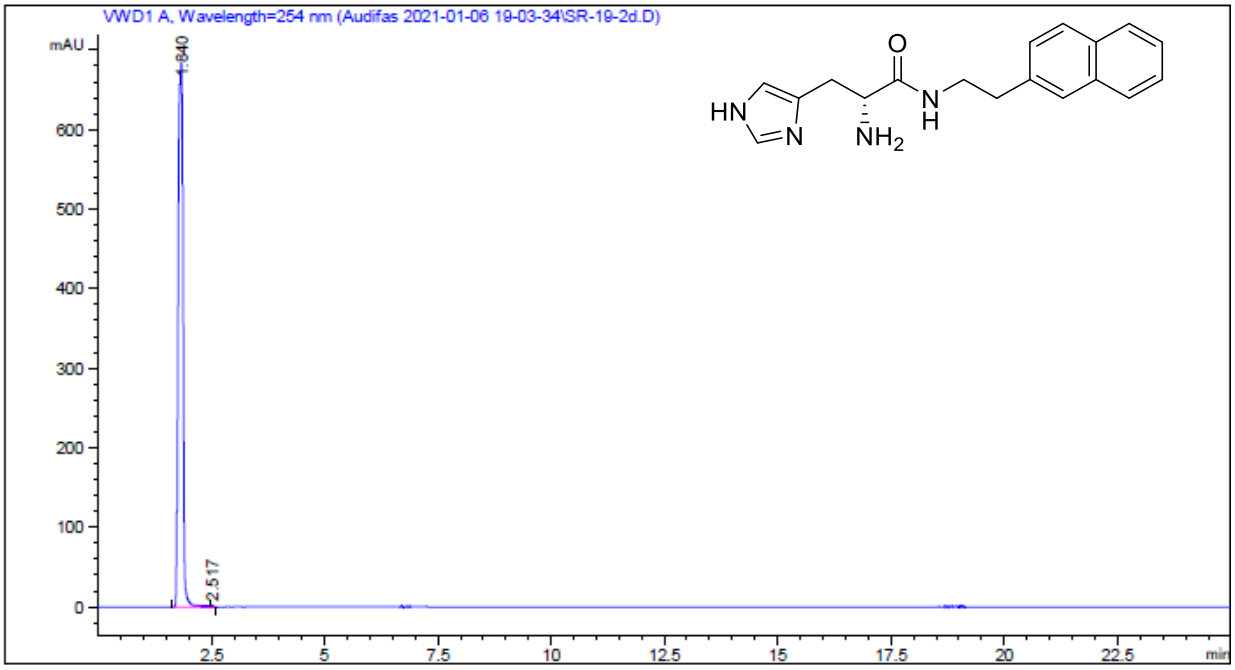
Totals :                  5478.52557  714.02272
```

(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(2-(naphthalen-2-yl)ethyl)propanamide **(9g)**

```
Sample Name: SR-19-2d

======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   5
Acq. Instrument : 1220 HPLC                       Location :   63
Injection Date  : 1/6/2021 8:59:17 PM                 Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR-19-2d
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.840 BV R  0.1194 5018.99854  685.26447  99.8394
   2   2.517 VB E  0.0575    8.07406    1.92153   0.1606

Totals :                  5027.07260  687.18600
```

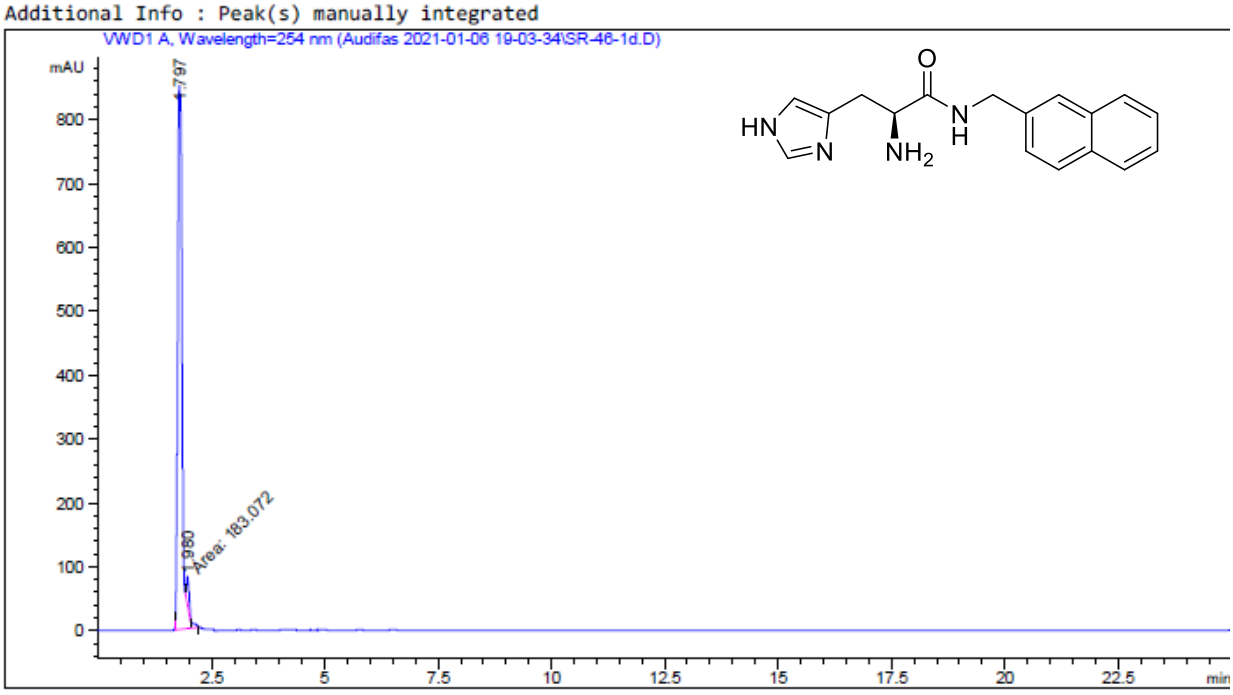
(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-2-ylmethyl)propanamide **(10a)**

```
Sample Name: SR-46-1d

==================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   4
Acq. Instrument : 1220 HPLC                     Location :   62
Injection Date  : 1/6/2021 8:31:47 PM               Inj :   1
                                               Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR-46-1d

Additional Info : Peak(s) manually integrated
```




```
==================================================================
                      Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.797 MM R  0.1145 6074.97803  852.00824  97.0746
   2   1.980 MM T  0.0675  183.07237   45.17344   2.9254

Totals :                  6258.05040  897.18168
```

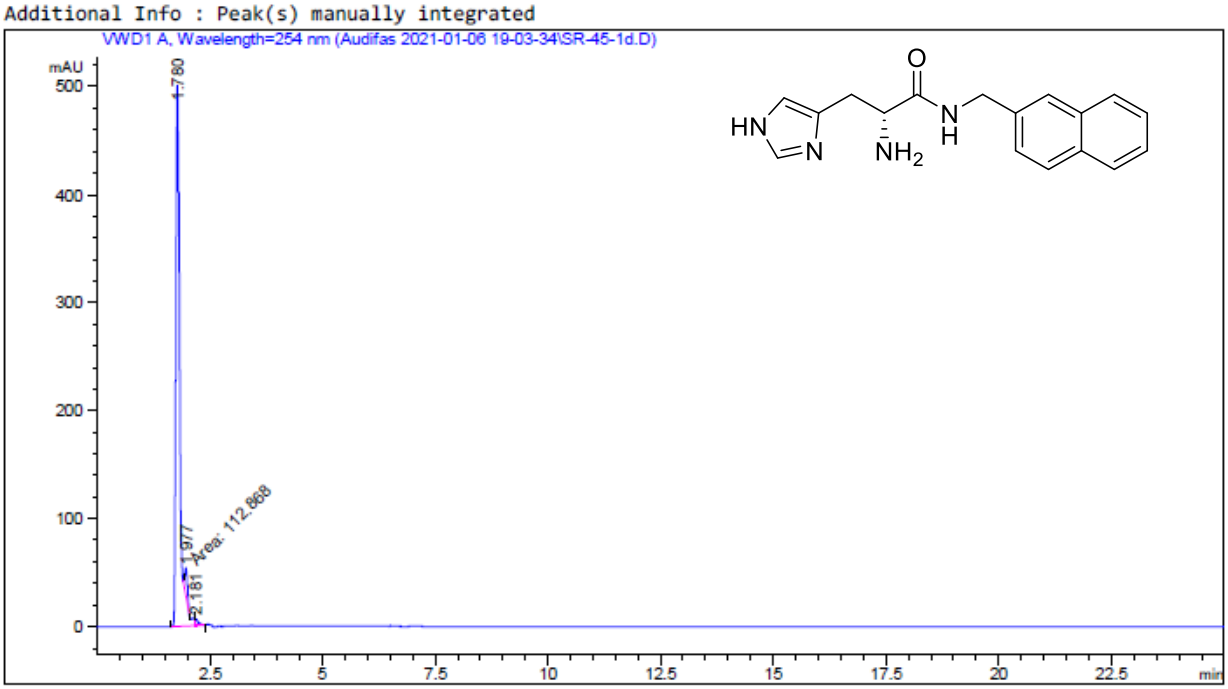
(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-2-ylmethyl)propanamide **(10b)**

```
Sample Name: SR-45-1d

==================================================================
Acq. Operator   : SYSTEM                       Seq. Line :   3
Acq. Instrument : 1220 HPLC                     Location :   61
Injection Date  : 1/6/2021 8:04:18 PM               Inj :   1
                                               Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water TFA 0.05 %

Sample Info     : SR-45-1d

Additional Info : Peak(s) manually integrated
```




```
==================================================================
                      Area Percent Report
==================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

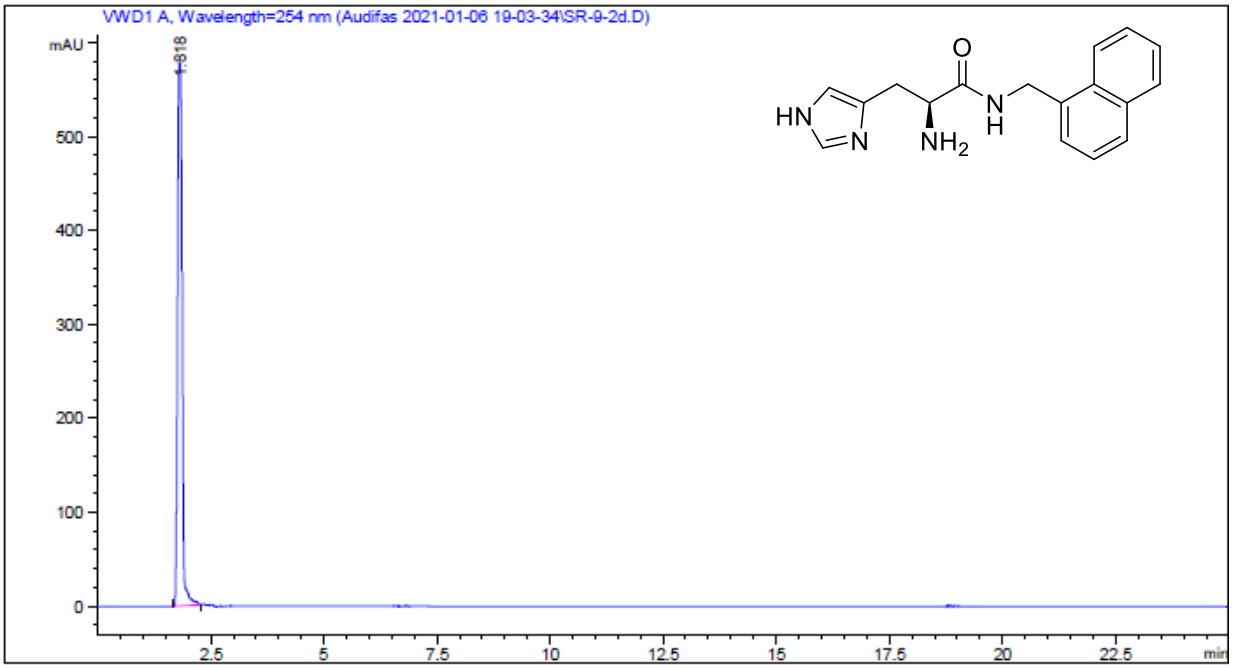
Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.780 BV R  0.0867 3096.73462  500.99149  95.4401
   2   1.977 MM T  0.0721  112.86839   26.09600   3.4786
   3   2.181 VB E  0.0706   35.08551    6.10260   1.0813
```

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-1-ylmethyl)propanamide **(10c)**

```
Sample Name: SR-9-2d

==========================================================================
Acq. Operator   : SYSTEM                          Seq. Line :   7
Acq. Instrument : 1220 HPLC                       Location :   65
Injection Date  : 1/6/2021 9:54:18 PM                  Inj :   1
                                                 Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water  TFA 0.05 %

Sample Info     : SR-9-2d
```




```
==========================================================================
                        Area Percent Report
==========================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.818 BV R  0.1146 4119.67236  578.71716 100.0000

Totals :                  4119.67236  578.71716
```

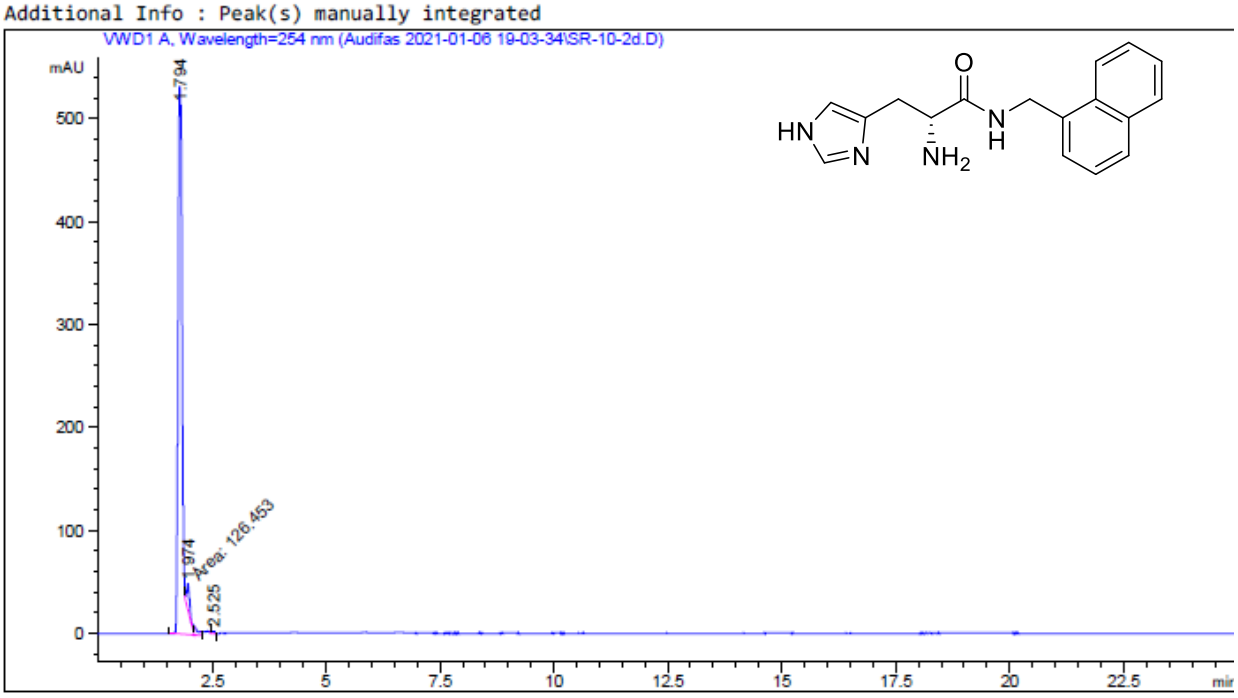
(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(naphthalen-1-ylmethyl)propanamide **(10d)**

```
Sample Name: SR-10-2d

==================================================================
Acq. Operator   : SYSTEM                          Seq. Line :  10
Acq. Instrument : 1220 HPLC                        Location :  66
Injection Date  : 1/6/2021 11:17:05 PM                 Inj :  1
                                                  Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 30.000 µl
Method          : C:\Chem32\1\Data\Audifas 2021-01-06 19-03-34\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 1/6/2021 7:13:21 PM by SYSTEM
Method Info     : Methanol/Water amonnia 0.05%

Sample Info     : SR-10-2d

Additional Info : Peak(s) manually integrated
```




```
==================================================================
                      Area Percent Report
==================================================================

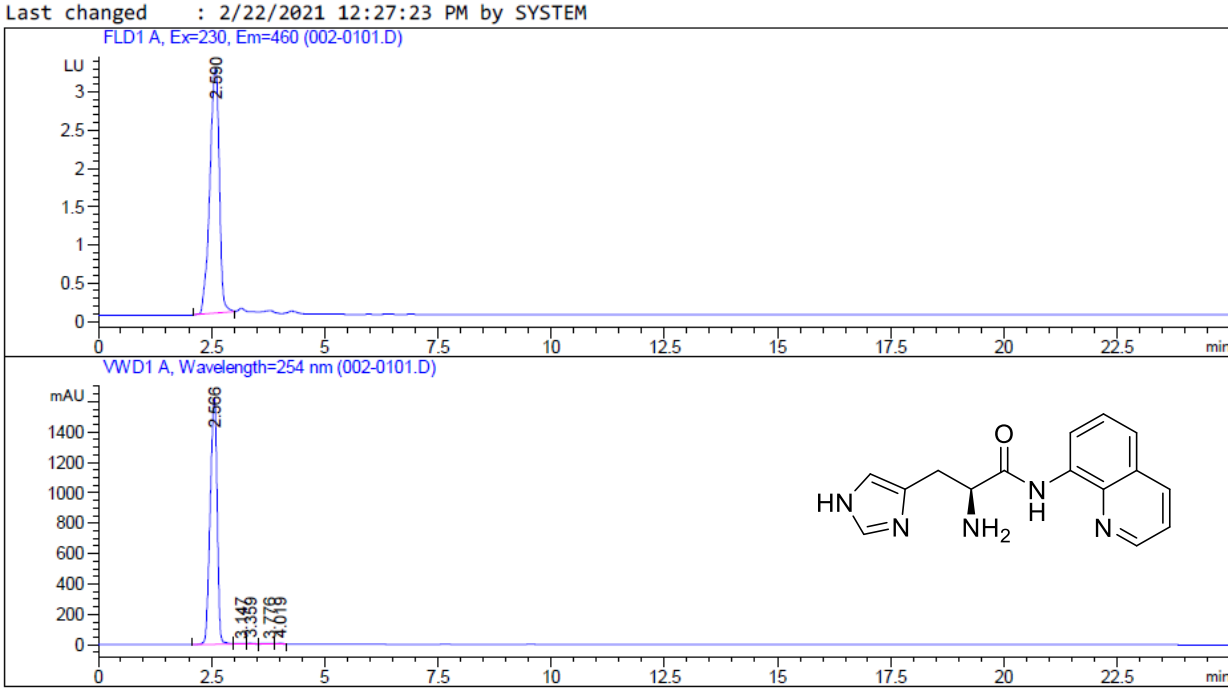
Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   1.794 MM R  0.1115 3561.12109  532.48346  96.2611
   2   1.974 MM T  0.0849  126.45335   25.51042   3.4182
   3   2.525 VB    0.0615   11.86573    2.30843   0.3207
```

(*S*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(quinolin-8-yl)propanamide **(11a)**

```
Acq. Operator   : SYSTEM                          Seq. Line :   1
Acq. Instrument : HPLC                           Location : Vial 2
Injection Date  : 2/22/2021 12:28:32 PM               Inj :   1
                                                 Inj Volume : 20.000 µl
Sequence File   : C:\Chem32\1\DATA\Ao\KK 2021-02-22 12-27-23\KK.S
Method          : C:\CHEM32\1\DATA\AO\KK 2021-02-22 12-27-23\CQ.M (Sequence Method)
Last changed    : 2/22/2021 12:27:23 PM by SYSTEM
```




```
=====================================================================
                    Area Percent Report
=====================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Do not use Multiplier & Dilution Factor with ISTDs


Signal 2: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.566 BV    0.1634 1.72670e4  1618.76880  99.0040
   2   3.147 VV    0.1739   57.22061    4.48363   0.3281
   3   3.359 VB    0.1301   38.78291    4.30338   0.2224
   4   3.776 BV    0.2168   30.98909    2.18032   0.1777
   5   4.019 VV    0.1189   46.72070    5.86617   0.2679

Totals :                 1.74407e4  1635.60230
```

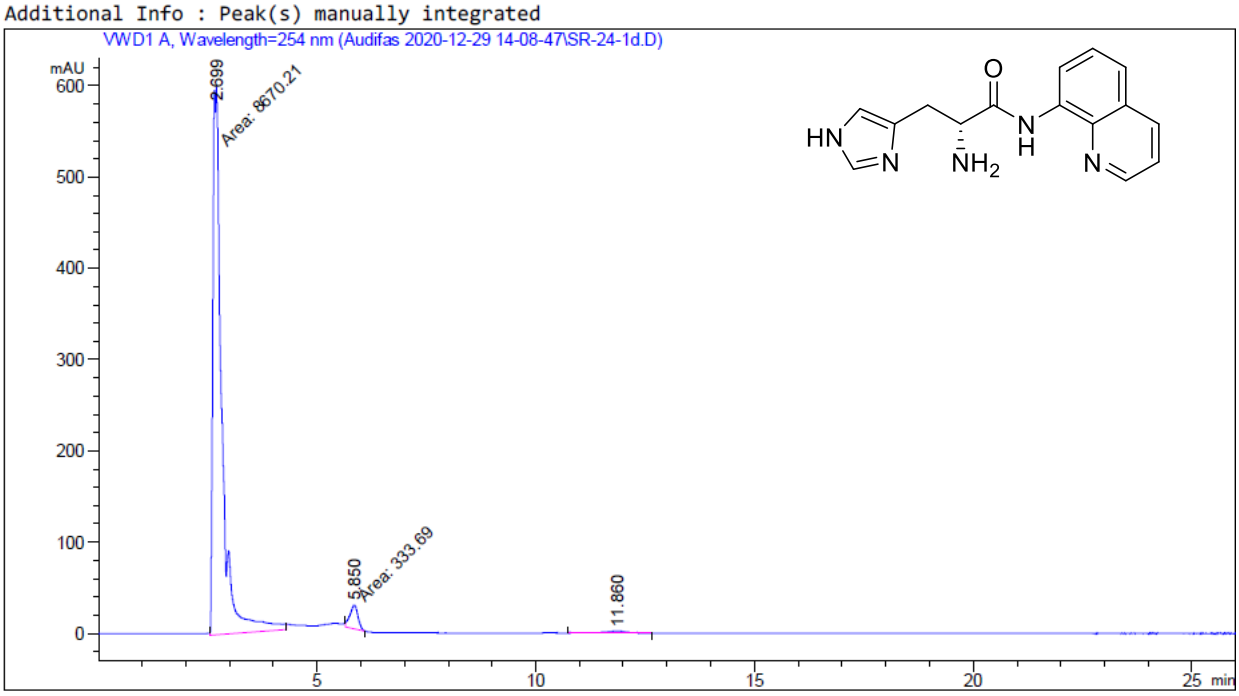
(*R*)-2-amino-3-(1*H*-imidazol-4-yl)-*N*-(quinolin-8-yl)propanamide **(11b)**

Sample Name: SR-24-1d

```
======================================================================
Acq. Operator   : SYSTEM                          Seq. Line :  17
Acq. Instrument : 1220 HPLC                        Location :  55
Injection Date  : 12/29/2020 9:52:27 PM               Inj :  1
                                                  Inj Volume : 50.000 µl
Different Inj Volume from Sample Entry!  Actual Inj Volume : 25.000 µl
Method          : C:\Chem32\1\Data\Audifas 2020-12-29 14-08-47\Wash_Shafikur.M (Sequence
                  Method)
Last changed    : 12/29/2020 2:08:51 PM by SYSTEM
Method Info     : Wash_Shafikur    Water 0.05 %TFA / Methanol

Sample Info     : SR-24-1d

Additional Info : Peak(s) manually integrated
```




```
======================================================================
                        Area Percent Report
======================================================================

Sorted By             :      Signal
Multiplier            :      1.0000
Dilution              :      1.0000
Use Multiplier & Dilution Factor with ISTDs


Signal 1: VWD1 A, Wavelength=254 nm

Peak RetTime Type  Width     Area      Height     Area
  #   [min]        [min]   [mAU*s]     [mAU]       %
----|-------|----|-------|----------|----------|--------|
   1   2.699 MM T  0.2403 8670.20898  601.43866  95.4375
   2   5.850 MM T  0.2157  333.68973   25.78662   3.6731
   3  11.860 VB R  0.4245   80.79691    2.22593   0.8894
```